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# FLUSH: A Tool for the Design of Slush Hydrogen Flow Systems

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# FLUSH: A TOOL FOR THE DESIGN OF SLUSH HYDROGEN FLOW SYSTEMS

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## SUMMARY

As part of the National Aerospace Plane Project an analytical model was developed to perform calculations for in-line transfer of solid-liquid mixtures of hydrogen. This code, called FLUSH, calculates pressure drop and solid fraction loss for the flow of slush hydrogen through pipe systems. The model solves the steady-state, one-dimensional equation of energy to obtain slush loss estimates. A description of the code is provided as well as a guide for users of the program. Preliminary results are also presented showing the anticipated degradation of slush hydrogen solid content for various piping systems.

## INTRODUCTION

Solid-liquid mixtures of hydrogen are currently being considered for use in the National Aerospace Plane (NASP) Project. These mixtures, known as slush hydrogen, offer the advantages of greater density and heat capacity than normal-boiling-point liquid hydrogen. An important consideration in the use of solid-liquid mixtures of hydrogen is the ability to transfer the fluid through flow lines. Previous studies (refs. 1 and 2) have shown that slush hydrogen with a solid fraction of 0.5 flows well through lines, valves, and other flow restrictions. However, few studies have shown calculations of the expected solid loss due to melting in these flow systems. Ewart and Dergance (ref. 3) performed such a study for the use of densified cryogenic propellants on advanced space transportation systems, but this effort concentrated on line sizes greater than 6 in. in diameter and the code it produced is not available in the open literature. Therefore, as part of the NASP project, an analytical model that performs calculations on the flow of slush in vacuum-jacketed lines was developed. This code, called FLUSH (FLOW of SLUSH), calculates the pressure drop and slush hydrogen solid fraction loss for steady-state, one-dimensional flow. These calculations can be useful in determining slush transfer parameters for experimental line analyses as well as parameters important in the design of slush transfer lines to the NASP vehicle itself.

This report summarizes the method of solution used in the code and provides a guide to the users of the code. The program solves the steady-state, one-dimensional energy equation and the Bernoulli equation for pipe flow. Slush hydrogen and liquid hydrogen properties are obtained by using the GASPLUS (ref. 4) physical properties code. The users guide gives input parameters required for FLUSH and example problems on the use of the code. The input requirements of the program include flow rate, upstream pressure, initial solid hydrogen fraction, element heat leak, and element parameters such as length and diameter. FLUSH also allows the user to specify pipe components such as elbows, valves, and straight pipe segments. FLUSH further allows the user to input friction factors and has the capability of calculating heat leak rates

for vacuum-jacketed piping, if estimates of the heat leak are not available. If standard insulation is desired instead of vacuum-jacketed piping, the user may specify input parameters that will allow the program to calculate heat leak for these cases. In addition, the user may vary the flow rate to obtain the hydrogen solid fraction loss for a range of flows. A FORTRAN program listing of the code is provided, and tables that give basic engineering data for designing vacuum-jacketed piping systems are presented.

In addition to the method of solution and the users guide, a summary of initial results obtained with the code is presented. The solid fraction degradation calculated by FLUSH is compared with estimates from previous analyses. Tradeoffs between heat leak and friction losses are discussed for various pipe sizes and lengths. Finally, a preliminary calculation of the slush loss expected at the NASA Lewis Research Center Plum Brook Station's slush hydrogen test facility in Sandusky, Ohio, is presented.

### FLUSH METHOD OF SOLUTION

The following section describes the method of solution used to calculate pressure drop and flow rate for slush hydrogen piping systems. Appendix A defines the symbols used below. Appendix B gives the actual program listing for FLUSH.

FLUSH solves the steady-state, one-dimensional energy equation given by

$$\Delta H + \frac{\Delta V^2}{2g_c J} + \frac{q}{g_c J} \Delta Z = \frac{Q}{\dot{m}} - W_s \quad (1)$$

and the Bernoulli equation given by

$$\frac{144\Delta P}{\rho} + \frac{\Delta V^2}{2g_c} + \frac{q}{g_c} \Delta Z + F + JW_s = 0 \quad (2)$$

where

$$F = \frac{fV^2 L}{2g_c D} \quad (3)$$

In the analysis the following assumptions are made:

1. Steady-state flow exists in the system.
2. The fluid is incompressible.
3. The lines are precooled to triple-point temperature (assumed to be a constant) before the slush flows through the lines.
4. The slush is well mixed before entering the line, and the effects of settling are neglected.

5. The viscosity of slush is the viscosity of liquid hydrogen because the viscosity of slush hydrogen is not known.

6. Heat of fusion is a constant.

7. The heat leak into the slush is used to melt the solid until all the solid is gone; the temperature is assumed to stay constant during the melting process. Once all the solid has melted, the heat leak produces a temperature rise in the remaining liquid.

The assumptions that the fluid is well mixed and that the lines are pre-cooled are probably the most significant of those listed.

FLUSH solves the preceding equations for each element specified by the user. An element can be a straight pipe length, a valve, a fitting, or any other piece of equipment in a piping system. Each element has two nodes; one node is at the beginning of the element, described by a set of initial or entrance conditions, and the other node is at the end of the element and is characterized by the final or exit conditions.

The following steps are used in solving these equations for the pressure drop through an element:

1. Initialize temperatures, pressures, slush solid fraction, and heat leak, as well as element length, diameter, and height.

2. Use GASPLUS to obtain initial-element liquid hydrogen properties.

$$\rho_l = \rho_l(T_l, P_l) \quad (4)$$

$$\mu_l = \mu_l(T_l, P_l) \quad (5)$$

3. Calculate an effective slush density at the inlet.

$$\rho_{mix} = x_l \rho_s + (1 - x_l) \rho_l \quad (6)$$

4. Calculate mass flow rate and Reynolds number. The Reynolds number for slush is calculated from the liquid viscosities.

$$\dot{m} = \frac{60 \dot{q} \rho_{mix}}{7.48} \quad (7)$$

$$Re = \frac{4 \dot{m}}{3600 \pi \mu_l D} \quad (8)$$

5. Calculate a friction factor by using either standard liquid friction factor correlations (the default selection) or National Institute of Standards and Technology (NIST, formerly the National Bureau of Standards) slush correlations. The standard friction factor correlation used is the Colbrook equation, which is a fit of data from the Moody plot of friction factor versus Reynolds number.

6. Calculate the downstream pressure from Bernoulli's equation.

$$P_f = P_i - \frac{\rho_{mix}}{144} \left( \frac{\Delta V^2}{2g_c} + \frac{g}{g_c} \Delta Z + F + JW_s \right) \quad (9)$$

Following calculation of the pressure drop the slush solid fraction exiting the element can be calculated. This is done in the following steps:

1. Obtain final-element hydrogen properties from GASPLUS.

$$\rho_l = \rho_l(T_f, P_f) \quad (10)$$

$$\mu_l = \mu_l(T_f, P_f) \quad (11)$$

2. Obtain initial- and final-element liquid enthalpies from GASPLUS.

$$H_{liq,f} = H_{liq,f}(T_f, P_f) \quad (12)$$

$$H_{liq,i} = H_{liq,i}(T_i, P_i) \quad (13)$$

3. Calculate initial mixture enthalpy.

$$H_{mix,i} = H_{liq,i} - x_i H_{fus} \quad (14)$$

4. Calculate final mixture enthalpy from the energy equation.

$$H_{mix,f} = H_{mix,i} + \frac{Q}{\dot{m}} - \frac{\Delta V^2}{2g_c J} - \frac{g}{g_c J} \Delta Z - W_s \quad (15)$$

The heat leak into the pipe element  $Q$  is normally a constant provided by manufacturers of cryogenic equipment. The value of  $W_s$  is zero unless a pump is used in the system.

5. Calculate the change in slush solid fraction due to heat input.

$$x_{f,h} = \frac{H_{liq,f} - H_{mix,f}}{H_{fus}} \quad (16)$$

6. Calculate the change in slush solid fraction due to friction losses.

$$\Delta x_{frict} = \frac{F}{H_{fus} J} \quad (17)$$

7. Calculate the slush fraction at the element exit.

$$x_f = x_{f,h} - \Delta x_{frict} \quad (18)$$

If slush still exists at the exit, the calculations are repeated using the solid fraction  $x_f$  to obtain a new value of  $\rho_{mix}$ . The calculations are stopped when convergence is reached on density and pressure. The convergence criteria are 0.0001 lbm/ft<sup>3</sup> and 0.0001 psi, respectively.

If  $x_f = 0$ , calculations are performed to find the increase in liquid temperature due to heat input.

1. Obtain the initial element enthalpy, density, and viscosity by using GASPLUS.

2. Obtain the saturated liquid enthalpy and the saturation temperature.

3. Find the final liquid enthalpy from the energy equation.

4. Obtain a liquid temperature from GASPLUS.

$$T_{liq,f} = T_{liq,f}(H_{liq,f}, P_f) \quad (19)$$

If  $T_{liq,f} \geq T_{sat}$ , assume that  $T_{liq,f} = T_{sat}$  and make a comment in the output. The program does not perform calculations for liquid-gas two-phase flow.

5. Repeat calculations to obtain convergence on temperature.

Once the slush solid fraction or the temperature at the output of the element has been obtained, the calculations can be repeated for a new element. Following completion of the calculations on a system of elements the initial slush flow rate is increased, if the user chose to vary the flow rate, and the entire process is repeated.

The code provides several options for determining heat transfer into a pipe element. If industry estimates are available, a constant value is used. If these estimates are not available for a vacuum-jacketed pipe or if layers of insulation are used on the pipe element, the FLUSH code can calculate the heat transfer values.

If the user desires to estimate the heat transfer into a vacuum-jacketed pipe element, a subroutine is provided in the code to perform this calculation. The method of calculating this heat transfer rate is discussed by Barron (ref. 5) and is described here.

The primary source of heat transfer across a vacuum-jacketed pipe is radiation. A cross section of a vacuum-jacketed pipe is shown in figure 1. The heat rate can be obtained by the equation

$$Q = F_e F_{1-2} \sigma A_1 (T_2^4 - T_1^4) \quad (20)$$

where, for no radiation shields, the emissivity factor  $F_e$  is defined as

$$\frac{1}{F_e} = \frac{1}{\epsilon_1} + \frac{A_1}{A_2} \left[ \frac{1}{\epsilon_2} - 1 \right] \quad (21)$$

Many vacuum-jacketed pipes contain radiation shielding to reduce the radiant heat transfer. In this case the emissivity factor is defined as

$$\frac{1}{F_e} = \left[ \frac{1}{\epsilon_1} + \frac{1}{\epsilon_s} - 1 \right] + (N - 1) \left[ \frac{2}{\epsilon_s} - 1 \right] + \left[ \frac{1}{\epsilon_2} + \frac{1}{\epsilon_s} - 1 \right] \quad (22)$$

Ewart and Dergance (ref. 3) point out that actual values of the emissivity factor can be considerably higher (a factor of five or greater) than those calculated. Therefore, in this study the actual emissivity used will be five times

that calculated. For cryogenic vessels the inner vessel is normally completely enclosed by the outer vessel, and  $F_{1-2} = 1$ .

For most vacuum-jacketed lines the vacuum space pressure is  $1 \times 10^{-5}$  torr or less. In this case, as shown by Ludtke and Voth (ref. 6), molecular conduction has little effect on the heat transfer. Therefore, calculations for molecular conduction in the vacuum space have not been included in the subroutine.

Note that the industry estimates of heat leak and those calculated by the subroutine are steady-state heat leak values. In an actual transfer line the heat leak may be higher owing to chilldown effects. Although this code assumes the line is chilled before the slush is placed in the line, it is possible to calculate transient effects (ref. 5). This transient analysis, however, is beyond the scope of the present work.

For most ground applications of slush hydrogen, vacuum-jacketed piping would be used for transfer. There may be cases, however, where vacuum-jacketed line is not feasible. In this case standard insulations may be required. Figure 2 shows a crosssection of piping with insulation. For conduction through insulation the heat transfer is calculated as

$$Q = UA_O(T_{air} - T_{fluid}) \quad (23)$$

where

$$\frac{1}{UA_O} = \frac{1}{h_i A_i} + \frac{\Delta r_a}{k_a A_{a,lm}} + \frac{\Delta r_b}{k_b A_{b,lm}} + \frac{\Delta r_c}{k_c A_{c,lm}} + \frac{\Delta r_d}{k_d A_{d,lm}} + \frac{1}{h_o A_o} \quad (24)$$

where the subscripts a, b, c, and d refer to the pipe material and the various types of insulation.

The log-mean area is defined as

$$A_{lm} = 2\pi L(r_2 - r_1) / \left[ \ln\left(\frac{r_2}{r_1}\right) \right] \quad (25)$$

For the case where a pump exists in the flow lines, additional calculations are required. For a pump the one-dimensional energy equation and Bernoulli equation (eqs. (1) and (2)) are solved with the addition of a shaft work term as discussed previously. The pump shaft work is defined as

$$W_s = - \frac{60(HP)}{0.02356m} \quad (26)$$

In addition, the pressure rise can be calculated once the efficiency of the pump and the shaft work are known, as given in equation (2).

$$\frac{-144\Delta P}{\rho} = JW_{s,min} + \frac{\Delta V^2}{2g_c} + \frac{q}{g_c} \Delta Z \quad (27)$$

where  $W_{s,min}$  is the minimum shaft work required, or the hydraulic horsepower. This hydraulic horsepower is defined as

$$W_{s,min} = F/J + W_s = nW_s \quad (28)$$

## FLUSH USERS GUIDE

The standard input file for the FLUSH code is divided into two sections: a namelist section and an element description section. In addition, if it is desired to calculate the heat leak into any element, a radiation heat transfer parameter section must be added. This section is not required if the heat leak rates are known. Also, if standard insulation or bare pipes are used, two additional sections are required. These are described below. The input file must be named "FLUSH.INP" for the code to run. Additional notes for users on the NASA Lewis Research Center VAX system are provided in appendix C.

In the namelist \$PARAMS section the variables are defined as follows:

FLOMIN	minimum flow rate, if it is desired to vary the flow in the analysis, gpm
FLOMAX	maximum flow rate, gpm
FLOINC	flow increment, gpm
E	pipe roughness, ft
P	initial (upstream) pressure, psi
XI	initial slush solid fraction
ZIN	initial (entrance) height of element 1, ft
NUM	number of elements in system
OUT	output control flag. If $OUT \leq 0$ , only final output conditions will be printed to unit 9; otherwise, output conditions of each element are displayed
LONG	output control flag for individual variable output. If $LONG \geq 0$ , the individual variables are printed to unit 8. This is useful in examining properties, etc.
XPUT	input control flag. If $XPUT \geq 0$ , print input data
FRICT	NIST friction factor option. If $FRICT \geq 0$ , use standard friction factor correlations; otherwise, use NIST relations for slush
HEAT	element heat transfer flag. If $HEAT \geq 0$ , heat transfer into each element has units of Btu/hr ft; otherwise, units of Btu/hr are used
TAMB	ambient temperature, °R
VAC	number of elements that require a calculation for the heat leak rate
XRS	number of radiation shields in vacuum-jacketed line. This is required for the heat leak calculation. Note: XRS is equivalent to N in equation (22)



NOJAC     number of elements that do not have vacuum jacketing (see eqs. (23) and (24))

The element description section of the input requires the information necessary to define an element. These variables are defined as follows:

ELE       element number

TYPE      element type. The codes corresponding to the element type are given in table 1

DIA       element diameter, in.

LENGTH   element length, ft

TEMP      element temperature, °R

Q          element heat transfer rate, Btu/hr or Btu/ft hr

K          flow resistance coefficient, normally used to define valves, pipe bends, and fittings

HT        height of element at its exit, ft

Table 2 shows an example of a standard input file for the FLUSH program. In this example an analysis is performed on a 100-ft length of 1.5-in.-diameter Schedule 5S vacuum-jacketed pipe. The line is divided into five sections, with bellows and bayonet fittings installed in the line. In addition, two globe valves are placed in the line, one near the beginning of the flow and one near the end. The entire line is located 25 ft above the ground. It is desired to vary the flow from a minimum of 10 to 100 gpm in increments of 5 gpm and find how much slush is lost at the end of the flow length given an initial solid fraction of 0.5. In this case conditions at the end of the piping system will be printed for each flow rate; input conditions also will be printed. Standard friction factor correlations will be used, and the heat transfer will be defined in British thermal units per hour foot for each element. The ambient temperature is 530 °R, and no heat leak calculations are required (VAC = 0). The variable XRS is not needed here as no heat leak estimates are being calculated, but it is shown for completeness.

The element conditions are defined in the second section of the input. The temperature of the slush hydrogen is assumed to be 25.2 °R. This value for the triple-point temperature is required, rather than the normal value of 24.84 °R, owing to the limitations of GASPLUS. The heat transfer rates to each of the individual elements were obtained from estimates made in industry; therefore, the radiation parameter section is not required. Estimates of heat leak to liquid hydrogen for various size pipes are shown in table 3.

As discussed in the literature (ref. 7), valves are normally defined in terms of either a flow coefficient  $C_v$  or a resistance coefficient  $K$  where

$$C_v = \frac{29.9d^2}{\sqrt{K}} \quad (29)$$

$$K = f \frac{L}{D} \quad (30)$$

Typical  $C_v$  and heat leak values for valves and fittings are given in table 4. If  $K$  is set equal to zero, the code assumes that a straight vacuum-jacketed line is being used.

Output from this example is shown in table 5. The listing gives the initial conditions of each element in the first page of output. The output parameters include the initial volumetric flow rate in gallons per minute, the mass flow rate in pounds per second, the final solid fraction, the change in solid fraction, and the pressure drop in pounds per square inch. Had all the solid hydrogen melted before the piping system exit had been reached, the heat leakage through the lines would be used to increase the temperature of the liquid hydrogen remaining. In this case the output under the heading "Change in x" would show the final fluid temperature. If the variable OUT had been specified as greater than zero, the conditions at each node would have been listed in the output. An example of this output is given in table 6. In addition to the previously defined parameters this form of output includes the pressure at each node (a node is defined as a point at the exit or entrance of an element).

Table 7 shows an example input file for the case where the heat leak rate is calculated for various elements of the previous example. In this case it is desired to calculate the rate of heat leak into elements 1 and 9. The variable VAC is set equal to 2, and the heat leak rate  $Q$  is set less than zero for elements 1 and 9 in the second section of the input. In addition, radiation heat transfer parameters must be defined: the element number, the outside diameter (o.d.) of the inner pipe, the inside diameter (i.d.) of the outer pipe, the emissivity of the inner pipe  $\epsilon_1$  ( $E_i$  in the input), and the emissivity of outer pipe,  $\epsilon_2$  ( $E_o$  in the input). The i.d. and o.d. for Schedule 5S stainless steel pipe are given in table 8. For this example a 3-in. pipe is used for the outer pipe, and 10 radiation shields between the inner and outer pipes are used as defined by  $XRS = 10$ . Note that the user must leave a space between the element description section and the radiation heat transfer parameter section. Also, the heat leak must be defined in British thermal units per hour foot when calculating the heat leak. Typical output for this example is given in table 9.

The output for this example shows a calculated heat leak of 0.552 Btu/hr ft, a value somewhat lower than the estimates given in table 3. These differences could be the result of differences in the surrounding conditions assumed for the industry estimates. Figure 3 shows the calculated heat leak as a function of ambient temperature for various pipe sizes. These calculations were based on vacuum-jacketed line with 10 radiation shields. Emissivities were assumed to be 0.3 for both the inner and outer lines. This emissivity value corresponds to that of polished stainless steel (ref. 8). Ten radiation shields with emissivities of 0.03 were assumed for the analysis. Comparing the heat leak values given in the figure with industry estimates of heat leak shows that the calculated values are generally between 5 and 13 percent lower than industry estimates, assuming an ambient temperature of 540 °R. Therefore, the calculated values of heat leak are expected to be close to those given by industry as design values.

When standard insulation is used, the type of insulation and the thickness must be specified. An example of such a case is shown in table 10. Here, in the first part of the insulation parameters, ELE is the element number, O.D. is

the pipe outer diameter, and T1, T2, and T3 are the thicknesses of the insulation. Up to three insulation layers are allowed by the code. In the second section of the insulation parameters the material types are specified. The material types corresponding to the codes are given in table 11. In this example five elements have either standard insulation or bare pipes, as specified by NOJAC=5. Various combinations of pipe materials and insulation schemes are shown. Note that the value of Q given in the element description section will be replaced by that calculated in this section; therefore, the Q specified will not be that given in the output. The output for this example is shown in table 12.

Finally, an example is presented in table 13 for the case where a pump exists in the flow system. In order to specify a pump in the line, the element type is set as either PP or PH. Element type PP signifies a pump with pressure rise and efficiency specified, while PH represents a pump with brake horsepower and efficiency given. In the example shown element 2 is set to type PH. In this case the horsepower is defined in the column under Q and the efficiency is input in the column under K. Also, element 4 is specified as type PP. The pressure drop for the pump in element 4 is input under the Q column and efficiency is placed in the column under K. Table 14 shows the output for the example; slush loss would occur in the pump owing to energy input by the pump.

The user should be aware that there are some limits on the code. These limits are summarized as follows:

1. The triple-point temperature of slush must be set equal to 25.2 °R, rather than 24.8 °R, owing to limitations on GASPLUS.
2. The exit pressure on the piping system must be greater than 1.1 psi (approximately the triple-point pressure of hydrogen). If this limit is not met because the inlet pressure was chosen to be too low, a message appears in the output file.
3. The maximum number of pipe elements the user can input is 90.
4. When all the slush has melted and the heat leak into the system causes a rise in temperature, the code is limited to a final temperature equal to that of the normal-boiling-point saturated liquid; gas-liquid two-phase flow is not included in FLUSH.
5. When using the NIST friction factor correlations, a maximum Reynolds number of only  $1 \times 10^6$  is possible because of limits on the data.
6. Up to three layers of insulation can be used for the pipe if vacuum-jacketed piping is not used.

## RESULTS

In order to obtain confidence in the code, a comparison was made with results from a previous analysis performed by Ewart and Dergance (ref. 3). In this study propellant heating due to heat leak and friction losses was considered, as in the present study. The losses due to friction were assumed to be the same as those for liquid hydrogen in the Ewart and Dergance study; therefore, standard friction factor correlations were used in this comparison. The

line considered was a 12-in.-diameter Schedule 5S pipe, 1750 ft long, with a constant heat leak of 12 Btu/hr ft. This heat leak value was calculated for a vacuum-jacketed pipe with multilayer insulation.

Figure 4 compares the slush loss calculated by FLUSH with that of the Ewart and Dergance study. As can be seen from the data, the results are nearly the same over the entire range of flow rates. Also, as discussed in the previous study, a minimum slush loss is reached at a flow rate between 3000 and 4000 gpm owing to tradeoffs in environmental heating and friction heating. At lower flow rates heat leak into the line is the major source of slush loss. At higher flow rates friction heating becomes significant. For example, in the Ewart and Dergance comparison at 500 gpm, the solid fraction loss due to environmental heating is 0.043, while that due to friction losses is essentially zero. At 10 000 gpm, the solid fraction loss due to environmental heating is 0.013, while the loss due to friction is 0.012. For cases with less heat leak the slush loss due to friction becomes the dominant loss mechanism.

Following a comparison against previous results the FLUSH code was run using various line sizes and pipe lengths. The results of these runs are shown in figure 5 for line lengths of 100, 200 and 500 ft, respectively. Each line was assumed to be only a straight section of Schedule 5S pipe, with no valves or fittings. As seen in each of these figures, a minimum solids loss was reached as in the case discussed previously. As the line size is decreased, it appears that the minimum occurs at a lower flow rate and at a higher solid fraction loss. This is due to the increased friction loss for smaller pipes, as can be seen in the equation

$$F = \frac{fV^2L}{2g_cD} \quad (3)$$

where  $F$  is the energy loss due to friction. The effect of a smaller pipe cross-sectional area is to increase the velocity which, coupled with a decreased diameter, leads to an increased friction loss. Hence, for a small pipe size the friction loss will begin manifesting itself at a lower flow rate, and, as the flow rate increases, the smaller pipe size will present higher solid fraction losses than do the larger pipe sizes. At low flow rates the larger pipes exhibit large slush losses. This is due to the higher surface area and hence higher heat leak into the larger pipes. For longer pipes the solid fraction losses increase owing to increases in both the friction loss and the environmental loss.

A piping system is presently being designed for the transfer of slush hydrogen at NASA Lewis Research Center's Plum Brook Station (ref. 9). This system, located in the K-Site facility, will require the transfer of slush hydrogen from a slush generator located approximately 100 ft from the slush test tank through 1.5-in.-diameter Schedule 5S pipe. An analysis of this piping system was performed by using the FLUSH code. The current system design contains two valves and five 20-ft sections of pipe. Bayonet fittings are used to connect these lines, and bellows are included in each of the lines. The heat leak used for the piping was obtained from industry estimates, provided in tables 3 and 4. The heat leak to the bellows was assumed to be that of the flexline. The roughness of the pipe was assumed to be 0.00015, typically the roughness of commercial steel. The flow rate was varied between 10 and 180 gpm and values of the slush loss were obtained.

The results of the slush loss analysis for K-Site are shown in figure 6. The results show that a fairly broad minimum occurs for this system. Between 35 and 60 gpm the solid loss is approximately 0.5 percentage point. In other words, the expected slush quality, if the fluid is transferred at a rate between 35 and 60 gpm, will be 49.5 percent when it reaches the test tank if the slush is initially 50 percent solid hydrogen. This represents a small loss when using standard vacuum-jacketed pipe with multilayer insulation. Therefore, helium transfer technology is not required for this system.

Finally, a comparison was done between the K-Site system for standard friction factor correlations and the same system for NIST friction factor results (refs. 1, 2, and 10). The experimental NIST friction factor results, shown in figure 7, were obtained on a transfer line 0.652 in. i.d., and 48.4 ft long. The data were obtained for Reynolds numbers less than  $8 \times 10^5$  and extrapolated to the lower flow rates. The solid line in the figure represents a standard friction factor correlation used by NIST. Because these friction factors were obtained empirically, it is not known how the friction factors will change outside the range of Reynolds numbers tested. In addition, the NIST results were obtained for smooth pipes. Therefore, the K-Site case discussed above was run again assuming a smooth pipe. The comparison is shown in figure 8 for an initial slush solid fraction of 0.5. As expected, the slush loss calculations are nearly identical at flow rates less than the minimum value because environmental heating dominates the loss of slush. At higher flow rates the slush loss estimates are quite close owing to the similar values of the friction factor. For the case using the NIST correlations values of slush loss are not shown for flow rates greater than 175 gpm owing to the constraint that the Reynolds number must be less than  $1 \times 10^6$ .

An additional issue that has not been addressed is that of critical velocity, the velocity where the solid hydrogen particles begin to settle. A previous study (ref. 1) has shown that for a 0.652-in.-i.d. pipe the critical velocity is 1.5 ft/sec. If this value is applied to larger pipe sizes, the critical flow rates range from 5.2 gpm for a 1-in.-diameter Schedule 5S pipe to 568 gpm for a 12-in.-diameter pipe, as shown in table 15. Below the critical flow rate the fluid becomes stratified, and the assumptions of one-dimensional flow and a well-mixed fluid are no longer valid. Therefore, the values produced by the code are not expected to be accurate at flow rates below the critical value. It is clear that more testing is required to obtain the critical velocities at larger pipe sizes as the values in table 15 are extrapolated from data for small lines. It has been postulated that the critical velocity depends on pipe diameter with the dependence ranging from direct proportionality to diameter to the 0.15 power (ref. 11).

#### CONCLUDING REMARKS

An analytical model was developed as part of the National Aerospace Plane Project to predict pressure drop and hydrogen solid fraction loss for a range of pipe sizes and flow rates. This code, called FLUSH, has direct application to the design of piping systems for slush hydrogen experimental test systems and for the flight and ground handling systems of the NASP vehicle. The code solves the one-dimensional, steady-state heat transfer equation to find the solid fraction loss. A listing of the program is provided, and a guide to the code's use is presented.

Results using FLUSH show that this code agrees well with previous analyses. Also, the data show that at low flow rates heat leak effects dominate the solid fraction loss, while at higher flow rates friction heating causes most of the solid hydrogen degradation. These two effects result in a minimum slush solid fraction loss for each pipe size. Further, as the pipe diameter increases, this minimum slush solid fraction loss decreases owing to the decreasing friction losses. An analysis of the NASA Plum Brook slush facility transfer line showed small losses for the piping system design being considered. The analysis will be compared with the test results upon completion of the facility. However, solid hydrogen losses appear to be in an acceptable range for fully chilled vacuum-jacketed lines designed for liquid hydrogen technology. Finally, comparing slush loss with standard friction factor correlations and with friction factor relations obtained in previous experiments showed similar values over the range of flow rates in this study.

# APPENDIX A

## SYMBOLS

$A_{a,lm}$	} log mean area of material a, b, c, or d, $ft^2$
$A_{b,lm}$	
$A_{c,lm}$	
$A_{d,lm}$	
$A_i$	inside surface area of pipe, $ft^2$
$A_o$	outside surface area of pipe, $ft^2$
$A_1$	area of surface 1 per foot, $ft^2/ft$
$A_2$	area of surface 2 per foot, $ft^2/ft$
$C_v$	flow coefficient
$D$	pipe diameter, ft
$d$	pipe diameter, in.
$F$	energy loss due to friction, $ft\ lbf/lbm$
$F_e$	emissivity factor
$F_{1-2}$	configuration factor
$f$	friction factor
$g$	acceleration due to gravity, $32.2\ ft/sec^2$
$g_c$	$32.2\ lbm\ ft/lbf\ sec^2$
$\Delta H$	change in enthalpy, $Btu/lb$
$HP$	brake horsepower, hp
$H_{fus}$	latent heat of fusion, $Btu/lb$
$H_{liq,f}$	enthalpy of liquid exiting element, $Btu/lb$
$H_{liq,i}$	enthalpy of liquid entering element, $Btu/lb$
$H_{mix,f}$	enthalpy of solid-liquid mixture exiting element, $Btu/lb$
$H_{mix,i}$	enthalpy of solid-liquid mixture exiting element, $Btu/lb$
$h_i$	inside heat transfer coefficient, $Btu/hr\ ft^2\ ^\circ R$
$h_o$	outside heat transfer coefficient, $Btu/hr\ ft^2\ ^\circ R$

J	mechanical equivalent of heat, 778.2 ft-lbf/Btu
K	resistance coefficient
$k_a$	thermal conductivity of material a, b, c, or d, Btu/hr ft R
$k_b$	
$k_c$	
$k_d$	
L	element length, ft
$\dot{m}$	mass flow rate, lb/hr
N	number of radiation shields
$\Delta P$	change in pressure, psi
$P_f$	pressure at element exit, psi
$P_i$	pressure at element entrance, psi
Q	heat leak into element, Btu/hr
q	volumetric flow rate, gpm
Re	Reynolds number (also R)
$\Delta r_a$	thickness of material a ( $r_a - r_i$ ), ft
$\Delta r_b$	thickness of material b ( $r_b - r_a$ ), ft
$\Delta r_c$	thickness of material c ( $r_c - r_b$ ), ft
$\Delta r_d$	thickness of material d ( $r_d - r_c$ ), ft
$r_a, r_b$	outside radius of material a, b, c, or d, ft
$r_c, r_d$	
$r_1$	inside radius for log-mean radius calculation, ft
$r_2$	outside radius for log-mean radius calculation, ft
$T_{air}$	air temperature, °R
$T_f$	temperature at element exit, °R
$T_{fluid}$	fluid temperature, °R
$T_i$	temperature at element entrance, °R



$T_{liq,f}$	liquid temperature at element exit, °R
$T_{sat}$	saturated liquid temperature, °R
$T_1$	temperature of surface 1 (equal to fluid temperature), °R
$T_2$	ambient temperature, °R
$U$	overall heat transfer coefficient, Btu/hr ft <sup>2</sup> °R
$V$	fluid velocity, ft/sec
$W_s$	shaft work, Btu/lb
$W_{s,min}$	minimum shaft work (hydraulic horsepower), Btu/lb
$x_f$	solid fraction at element exit
$x_{f,h}$	solid fraction at exit (heat leak considerations only)
$\Delta x_{frict}$	change in solid fraction due to friction losses
$x_i$	solid fraction at element inlet
$\Delta Z$	change in element height, ft
$\epsilon_s$	emissivity of radiation shielding
$\epsilon_1$	emissivity of surface 1
$\epsilon_2$	emissivity of surface 2
$\eta$	pump efficiency
$\mu$	triple-point hydrogen viscosity
$\mu_l$	liquid viscosity, lbm/ft sec
$\rho$	slush density, lbm/ft <sup>3</sup>
$\rho_{mix}$	solid-liquid mixture density, lbm/ft <sup>3</sup>
$\rho_s$	solid density, lbm/ft <sup>3</sup>
$\rho_l$	liquid density, lbm/ft <sup>3</sup>
$\sigma$	Stefan-boltzmann constant, $0.173 \times 10^{-8}$ Btu/hr ft <sup>2</sup> °R <sup>4</sup>

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APPENDIX B  
PROGRAM LISTING

# PROGRAM FLUSH

```

C
C-----
C      THIS PROGRAM CALCULATES PRESSURE LOSS AND SLUSH HYDROGEN
C      SOLID FRACTION LOSS IN STEADY STATE FLOWS IN PIPE SYSTEMS
C      (SIMPLE PUMP EQUATIONS ARE INCLUDED)
C
C      THE PROGRAM SOLVES THE STEADY-STATE, ONE DIMENSIONAL
C      ENERGY EQUATION.
C
C       $H2-H1 + (V2^{**2}-V1^{**2})/2GC + (G/GC)(Z2-Z1) = Q - WS$ 
C
C      AND THE BERNOULLI EQUATION
C
C       $(P2-P1)/RHO + (V2^{**2}-V1^{**2})/2GC + (G/GC)(Z2-Z1) + F + WS = 0$ 
C
C-----
C
C      TERRY HARDY
C      NASA LEWIS RESEARCH CENTER
C      CLEVELAND, OHIO 44135
C      (216) 433-2411
C
C      4-17-89
C
C-----
C
C      DATA ID/'PH2'/
C      DIMENSION EDIA(90),ELEN(90),ET(90),EX1(90),EZ2(90),EQIN(90),
C      ERHO(90),EP1(90),EMDOT(90),EQUAL(90),EDP2(90),
C      NODE(90),TS(90),HF1(90),ETNEW(90),ECLD(90),RE1(90),
C      NVAC(90),EVDI(90),EVD0(90),EEI(90),EE0(90),PTYP(90),
C      QFL(90),QL(90),PXTYP(10,90),INSUL(90)
C
C      COMMON /FLUX/ EP0D(90),ETI1(90),ETI2(90),ETI3(90),
C      NNOVJ(90),NOJAC,NXMAT(20,90),
C      EXLEN(90),EXDIA(90),EXT(90),TAMB,P,X1,FLOV,EPID(90),K
C
C      REAL MDOT,LEN,MPHR,LONG
C      CHARACTER*12 PXTYP
C      CHARACTER*10 TIM,TONE,TAM,TANE
C      CHARACTER*4 PTYP,TYPE,PTYP
C
C-----
C
C      READ IN INPUT DATA
C
C-----
C
C      OPEN(UNIT=7,FILE=' FLUSH.INP ', STATUS='OLD')
C
C      PARAMS NAMELIST SECTION
C
C      NAMELIST / PARAMS / FLOWIN,FLOMAX,FLOINC,E,P,X1,ZIN,NUM,OUT,
C      LONG,XPUT,FRICT,HEAT,TAMB,VAC,XRS,NOJAC
C      READ (7,PARAMS)
C      READ(7,800)
C      800)  FORMAT(A80)
C      NUMB = NUM
C      TOTLEN = 0.
C

```

**C**

31

```

C      XPUT = INPUT PRINT CONTROL FLAG
C      < 0 => DO NOT PRINT INPUT DATA
C      HEAT = HEAT TRANSFER PER ELEMENT FLAG
C      < 0 => BTU/HR, > 0 => BTU/HR-FT
C      FRICT = NBS SLUSH FRICTION FACTOR OPTION
C      < 0 => USE NBS CORRELATION FOR F
C      VAC = NUMBER OF ELEMENTS WHICH REQUIRE A CALCULATION FOR
C      THE HEAT LEAK RATE (VACUUM JACKETED PIPING)
C      XRS = NUMBER OF RADIATION SHIELDS FOR VJ LINE CALCULATIONS
C      TAMB = AMBIENT TEMPERATURE, R
C      NOJAC = NUMBER OF ELEMENTS WHICH REQUIRE A CALCULATION FOR
C      THE HEAT LEAK RATE (BARE PIPE OR STANDARD INSULATION)

```

```

C-----
C      CALCULATE THE HEAT LEAK RATE THROUGH VACUUM-JACKETED PIPES
C      IF VAC > 0
C

```

```

      IE = 1
      IF (VAC .LE. 0.0) GO TO 95
      DO 802 I=1,4
      READ(7,801)
801  FORMAT(A80)
802  CONTINUE
      DO 13 IB=1,NUMB
      NEWQ = EQIN(IB)
      TINRAD = ET(IB)
      IF (NEWQ .LT. 0.0) GO TO 14
      GO TO 13
14  READ(7,301) NELEM, DIN, DOUT, EI, EO
301  FORMAT(I4,2X,4F10.3)
      NVAC(IE) = NELEM
      EVDI(IE) = DIN
      EVDO(IE) = DOUT
      EEI(IE) = EI
      EEO(IE) = EO
      CALL VACUUM(QIN1,TAMB,XRS,DIN,DOUT,TINRAD,EI,EO)
      EQIN(IB) = QIN1
      IE = IE + 1
13  CONTINUE
95  CONTINUE

```

```

C-----
C      CALCULATE HEAT LEAK FOR PIPE WITHOUT VACUUM JACKET IF
C      NOJAC > 0
C-----
C

```

```

      IF (NOJAC .LE. 0) GO TO 150
      DO 311 I=1,4
      READ(7,310)
310  FORMAT(A80)
311  CONTINUE
      DO 213 I=1,NOJAC
      READ (7,312) NELE, POD, TI1,TI2,TI3
312  FORMAT(I4,4F10.2)
      N = NELE
      EPOD(N) = POD/12.0
      ETI1(N) = TI1
      ETI2(N) = TI2
      ETI3(N) = TI3

```

```

      NNOVJ(I)= NELE
213  CONTINUE
      DO 214 I=1,3
      READ(7,314)
314  FORMAT(A80)
214  CONTINUE
      DO 215 I=1,NOJAC
      READ(7,315) NELE, MAT1, MAT2, MAT3, MAT4
315  FORMAT(I4,6X,I5,5X,I5,5X,I5,5X,I5,5X)
      N      = NELE
      NXMAT(1,N) = MAT1
      NXMAT(2,N) = MAT2
      NXMAT(3,N) = MAT3
      NXMAT(4,N) = MAT4
      NNOVJ(I)  = NELE
      EXLEN(N)  = ELEN(N)
      EXDIA(N)  = EDIA(N)/12.0
      EPID(N)   = EDIA(N)/12.0
      EXT(N)    = ET(N)
215  CONTINUE
      FLOV = FLOMIN
      DO 216 I=1,NOJAC
      K = NNOVJ(I)
      CALL QFLUX(QFL(I),QL(I))
      IF (HEAT .LT. 0.) GO TO 217
      GO TO 218
217  EQIN(NNOVJ(I)) = QFL(I)
      GO TO 219
218  EQIN(NNOVJ(I)) = QL(I)
219  CONTINUE
216  CONTINUE
150  CONTINUE
C
C
      WRITE(9,500)
500  FORMAT(' FLUSH CODE'/)
C
C-----
C  CALL SYSTEM DATE AND TIME ROUTINES
C-----
C
      CALL DATE(TIM)
      CALL TIME(TONE)
      WRITE(9,130) TIM,TONE
130  FORMAT(/' ', A10,2X,A10)
C
C-----
C  WRITE INPUT PARAMETERS TO OUTPUT FILES IF XPUT . 0
C-----
C
      IF (XPUT .LT. 0.0) GO TO 42
      WRITE(9,405)
405  FORMAT(/'*****')
      WRITE(9,400)
400  FORMAT(/' INPUT PARAMETERS ')
      WRITE(9,401)
401  FORMAT(/' ELEMENT',2X,'TYPE',2X,'DIA.',1X,'IN.',3X,'LENGTH, FT',
      ,3X,'TEMP.',1X,'R',4X,'Q, BTU/HR-FT',3X,'K-F L/D '/')
      DO 43 IA=1,NUM

```

```

      IF (HEAT .LT. 0.0) GO TO 434
      GO TO 433
433    QALT = EQIN(IA)
      GO TO 435
434    QALT = EQIN(IA)/ELEN(IA)
435    CONTINUE
      WRITE(9,402) NODE(IA),PTYP(IA),EDIA(IA),ELEN(IA),ET(IA),
+ QALT,ECLD(IA)
402    FORMAT(I4,5X,A4,3X,F7.3,5X,F8.3,5X,F8.2,4X,F10.3,4X,F10.3)
      WRITE(9,403)
403    FORMAT('-----')
+ '-----')
43    CONTINUE
C
C      IF VAC > 0 PRINT RADIATION HEAT TRANSFER PARAMETERS
C
      IF (VAC .LE. 0) GO TO 97
      WRITE (9,410)
410    FORMAT(/' RADIATION HEAT TRANSFER PARAMETERS ')
      WRITE(9,406)
406    FORMAT(/' ELEMENT',4X,'INNER O.D.',2X,'OUTER I.D.',6X,
+ 'EI ',8X,'EO')
      DO 96 IC=1,VAC
      WRITE(9,407) NVAC(IC),EVDI(IC),EVDO(IC),EEI(IC),EEO(IC)
407    FORMAT(/I4,7X,F8.3,5X,F8.3,5X,F8.5,3X,F8.5)
96    CONTINUE
      WRITE(9,408)
408    FORMAT('-----')
+ '-----')
C
C      IF NOJAC > 0 PRINT BARE PIPE/STANDARD INSULATION PARAMETERS
C
97    IF (NOJAC .LE. 0) GO TO 197
      WRITE(9,411)
411    FORMAT(/' BARE PIPE OR STANDARD INSULATION PARAMETERS')
      WRITE(9,412)
412    FORMAT(/' ELEMENT',4X,'PIPE O.D.',3X,'INSULATION 1',
+ 4X,'INSULATION 2',4X,'INSULATION 3')
      WRITE(9,413)
413    FORMAT(25X,'THICKNESS, FT',3X,'THICKNESS, FT',3X,
+ 'THICKNESS, FT')
      DO 196, I=1,NOJAC
      K = NNOVJ(I)
      FORD = EPOD(K)*12.
      WRITE(9,414) K, FORD,ETI1(K),ETI2(K),ETI3(K)
414    FORMAT(/I4,10X,F7.4,6X,F7.4,9X,F7.4,9X,F7.4)
196    CONTINUE
      WRITE(9,419)
419    FORMAT(A80)
      WRITE(9,415)
415    FORMAT(/' ELEMENT',4X,'PIPE MATERIAL',4X,'INSULATION 1',
+ 4X,'INSULATION 2',4X,'INSULATION 3')
C
      DO 954 I = 1,NOJAC
      K = NNOVJ(I)
      IF (NXMAT(2,K) .LE. 0.0) GO TO 971
      IF (NXMAT(3,K) .LE. 0.0) GO TO 972
      IF (NXMAT(4,K) .LE. 0.0) GO TO 973
      GO TO 974
971    INSUL(I) = 1

```



```

GO TO 975
972 INSUL(I) = 2
GO TO 975
973 INSUL(I) = 3
GO TO 975
974 INSUL(I) = 4
975 CONTINUE
954 CONTINUE

```

C

```

DO 950 I = 1,NOJAC
DO 951 J = 1,4
K = NNOVJ(I)

```

C

C

C

C

C

MATCH MATERIAL CODE TO WORD DESCRIPTION

```

IF (NXMAT(J,K) .EQ. 101) GO TO 911
IF (NXMAT(J,K) .EQ. 102) GO TO 912
IF (NXMAT(J,K) .EQ. 103) GO TO 913
IF (NXMAT(J,K) .EQ. 104) GO TO 914
IF (NXMAT(J,K) .EQ. 201) GO TO 921
IF (NXMAT(J,K) .EQ. 202) GO TO 922
IF (NXMAT(J,K) .EQ. 203) GO TO 923
IF (NXMAT(J,K) .EQ. 204) GO TO 924
IF (NXMAT(J,K) .EQ. 205) GO TO 925
IF (NXMAT(J,K) .EQ. 301) GO TO 931
IF (NXMAT(J,K) .EQ. 302) GO TO 932
GO TO 942

```

```

911 PXTYP(J,K) = 'TEFLON'
GO TO 970
912 PXTYP(J,K) = '304 ST STEEL'
GO TO 970
913 PXTYP(J,K) = 'CARBON STEEL'
GO TO 970
914 PXTYP(J,K) = '6063 ALUM.'
GO TO 970
921 PXTYP(J,K) = 'FOAM GLASS'
GO TO 970
922 PXTYP(J,K) = 'SILICA'
GO TO 970
923 PXTYP(J,K) = 'POLYURETHANE'
GO TO 970
924 PXTYP(J,K) = 'POLYSTYRENE'
GO TO 970
925 PXTYP(J,K) = 'FOAM RUBBER'
GO TO 970
931 PXTYP(J,K) = 'FIBER GLASS'
GO TO 970
932 PXTYP(J,K) = 'PERLITE'
GO TO 970
942 PXTYP(J,K) = ' '
970 CONTINUE
951 CONTINUE
950 CONTINUE

```

```

DO 426 I=1,NOJAC

```

```

K = NNOVJ(I)

```

```

WRITE(9,427) K,PXTYP(1,K),PXTYP(2,K),PXTYP(3,K),PXTYP(4,K)

```

```

427 FORMAT(/I4,7X,A13,4X,A13,3X,A13,3X,A13)

```

```

426 CONTINUE

```





```

124  FORMAT(/' PIPE DIAMETER = ',F6.2,1X,'IN.',9X,'FLOW = ',
      +F7.2,2X,'GPM')
      WRITE(9,125) TOTLEN,XIN
125  FORMAT(/' TOTAL PIPE LENGTH = ',F8.2,1X,'FT',4X,'INITIAL
      + SOLID FRACTION = ',F5.3)
      WRITE(9,129) E, TO
129  FORMAT(/' PIPE ROUGHNESS = ',F9.6,9X,'INITIAL TEMPERATURE = ',
      +F7.2,1X,'R')
      WRITE(9,126)
      IF (XIN .LE. 0.0) GO TO 11
      WRITE(9,123)
123  FORMAT(/' NODE',3X,' PRESSURE',4X,
      +'FLOW, LB/S',5X,' FINAL SOLID',3X,' CHANGE IN X ',4X,
      +'PRESSURE, ')
      WRITE(9,128)
128  FORMAT(9X,'DROP,PSI',21X,'FRACTION',23X,' PSI'/)
      GO TO 12
C
11  WRITE(9,133)
133  FORMAT(/' NODE',3X,' PRESSURE',4X,
      +'FLOW, LB/S',5X,' FINAL SOLID',3X,' FINAL TEMP.,R ',3X,
      +'PRESSURE, ')
      WRITE(9,138)
138  FORMAT(9X,'DROP,PSI',21X,'FRACTION',23X,' PSI'/)
C
12  CONTINUE
126  FORMAT(/' .....
      + .....')
68  CONTINUE
C
C
C  REINITIALIZE VARIABLES
C
C
      HP      = 0.0
      WS      = 0.0
      EP1(1)  = PIN1
      EQUAL(1) = XIN
C
C
C  BEGIN CALCULATION LOOP FOR DIFFERENT PIPE ELEMENTS
C
C
      DO 32 J=1,NUM
      DIAM = EDIA(J)
      WS   = 0.0
      EFP  = 0.0
      LEN  = ELEN(J)
      T    = ET(J)
      TYPE = PTPA(J)
C
C
C  IF A PUMP IS USED, REINITIALIZE QIN AS HORSEPOWER (TYPE = PH)
C  OR PRESSURE RISE (TYPE = PP) AND CLD AS EFFICIENCY
C
C
      IF ((TYPE .EQ. ' PP') .OR. (TYPE .EQ. ' PH')) GO TO 781
      IF ((TYPE .EQ. ' PP ') .OR. (TYPE .EQ. ' PH ')) GO TO 781
      IF ((TYPE .EQ. 'PP ') .OR. (TYPE .EQ. 'PH ')) GO TO 781
      GO TO 782

```

```

782  QIN = EQIN(J)
      CONLD= ECLD(J)
      GO TO 783
781  EFFP = ECLD(J)
      IF (TYPE .EQ. ' PH') GO TO 785
      IF (TYPE .EQ. ' PH ') GO TO 785
      IF (TYPE .EQ. 'PH ') GO TO 785
      GO TO 784
785  HP = EQIN(J)
      GO TO 783
784  DPP = -EQIN(J)
783  CONTINUE
      P = EP1(J)
      X1 = EQUAL(J)
      Z2 = EZ2(J+1)
      Z1 = EZ2(J)
      ELEMI = NODE(J)
      JA = J + 1

C
C  SET CONSTANTS
C
      N = 0
      N1 = 0
      N2 = 0
      HFUS = 25.02
      TOLD = ET(J)
      PI = 3.14159
      GC = 32.2
      DIA = DIAM/12.0
      ED = E/DIA

      IF (TYPE .EQ. ' PP') .OR. (TYPE .EQ. ' PH')) GO TO 786
      IF ((TYPE .EQ. ' PP ') .OR. (TYPE .EQ. ' PH ')) GO TO 786
      IF ((TYPE .EQ. 'PP ') .OR. (TYPE .EQ. 'PH ')) GO TO 786

CHECK TO SEE IF HEAT LEAK INTO ELEMENT IS IN
BTU/HR OR BTU/FT-HR

      IF (HEAT .LT. 0.) GO TO 64
      GO TO 67
64  QINTOT = QIN
      GO TO 69
67  QINTOT = QIN*LEN
69  CONTINUE

C
786  CONTINUE

      RHOS = 5.40

      RHOS = DENSITY OF SOLID HYDROGEN, LB/CU.FT

CHECK TO MAKE SURE THE LAST NODE WAS NOT AT THE
SATURATION TEMPERATURE. IF SATURATION CONDITIONS EXIST
USE THE APPROPRIATE GASPLUS ROUTINES (NORMAL LIQUID
ROUTINES WILL NOT WORK FOR SATURATED CONDITIONS)

      IF (TS(1) .GE. ET(J)) GO TO 33
      GO TO 34
      RHOL = DLPSAT(P, ID)

```

```

      VIS1 = VLPSAT(P, ID)
      GO TO 35
34     CONTINUE
      RHOL = DENLPC(TOLD, P, ID)
      VIS1 = VISLPC(TOLD, P, ID)
35     CONTINUE
C
C     USE GASPLUS TO CALCULATE LIQUID HYDROGEN PROPERTIES
C     (VISCOSITY HAS UNITS OF LB/FT-SEC)
C
C-----
C     CALCULATE FLOW PARAMETERS
C-----
C
      RHOMIX = X1*RHOS + (1.-X1)*RHOL
      AREA   = PI*DIA**2*0.25
      EMDOT(J) = FLOW*RHOMIX*60./7.48
      MDOT    = EMDOT(1)
      VELI    = MDOT/(AREA*RHOMIX*3600.)
      RE      = DIA*RHOMIX*VELI/VIS1
      RE1(J)  = RE
C
C     CALL THE DESIRED FRICTION FACTOR ROUTINE
C
      IF (X1 .LE. 0.0) GO TO 17
      IF (FRICT .GE. 0.0) GO TO 17
      IF (RE .GT. 1.0E+06) GO TO 17
      GO TO 18
18     CALL FRICT1(X1, F, RE)
      GO TO 19
17     CALL FRICT2(F, RE, ED)
19     CONTINUE
C
C-----
C     SET EQUIVALENT PIPE LENGTH
C-----
C
C-----
C     IF A PUMP IS USED, THEN EFFICIENCY IS USED IN PLACE OF FLOW COEFF.
C-----
C
      IF (TYPE .EQ. ' PH') GO TO 788
      IF (TYPE .EQ. ' PH ') GO TO 788
      IF (TYPE .EQ. 'PH ') GO TO 788
      IF (TYPE .EQ. ' PP') GO TO 789
      IF (TYPE .EQ. ' PP ') GO TO 789
      IF (TYPE .EQ. 'PP ') GO TO 789
C
      IF (CONLD .GT. 0.0) GO TO 47
      GO TO 48
47     CNLD = CONLD/F
      GO TO 49
48     CNLD = LEN/DIA
49     CONTINUE
C
      GO TO 799
C
C-----
C     IF HORSEPOWER IS SPECIFIED CALCULATE SHAFT WORK, THEN PRESSURE
C-----

```

```

C
788  WS      = -HP*60./(.02356*MDOT)
      DP1A   = 0.0
      CNLD   = 0.0
      QINTOT = 0.0
      GO TO 790

```

```

C
C      WS IS IN UNITS OF BTU/LB
C

```

```

C      IF PRESSURE IS SPECIFIED, THEN CALCULATE MINIMUM SHAFT WORK, WSMIN,
C      ALSO KNOWN AS HYDRAULIC HORSEPOWER, THEN CALCULATE BRAKE
C      HORSEPOWER, HP
C

```

```

C
789  DP1B     DPP
      DP1A   = 0.0
      CNLD   = 0.0
      QINTOT = 0.0
      WSMIN  = DPP*MDOT*144./RHOMIX
      WSMIN IS IN UNITS OF FT-LB/HR
      HP1    = WSMIN/EFFP
      WS     = HP1/(MDOT*778.)
      HP     = HP1/(3600.*550.)
      GO TO 791

```

```

790  CONTINUE
      DP1A = (RHOMIX*CNLD*VELI**2*F)/(2.0*144.*GC)

```

```

791  CONTINUE
      DP1B = RHOMIX*(Z2-Z1)*(32.2/(GC*144.)) +
      QFLOW*778.*RHOMIX/144. + EFFP*WS*778.*RHOMIX/144.

```

```

792  CONTINUE
      DP1 = DP1A + DP1B
      IF ((ABS(Z2-Z1)-LEN) .GT. 0.00001) GO TO 610
      GO TO 620

```

```

793  WRITE(9,611) ELEMI
794  FORMAT(5X, ' CHANGE IN HEIGHT > ELEMENT LENGTH AT ELEMENT'.

```

```

795  F6.0)
      DELZ = ABS(Z2-Z1)
      WRITE(9,613) DELZ,LEN
796  FORMAT(5X, ' HEIGHT CHANGE = ',F7.2,1X,'FT.',2X,'LENGTH = ',
797  F7.2,'FT.')
```

```

798  WRITE(9,612)
799  FORMAT(8X, ' RE-EVALUATE SELECTED FLOW SYSTEM '/')
800  CONTINUE

```

```

-----
VELI  INITIAL FLOW VELOCITY, FT/SEC
RE    INITIAL FLUID REYNOLDS NUMBER
MDOT  FLOW RATE, LB/HR
F     FRICTION FACTOR
DP1   PRESSURE DROP FROM BERNOULLI'S EQUATION
-----

```

```

CALCULATE THE CHANGE IN SOLID FRACTION
-----

```

```

X2   X1

```

```

30  IF (X2 .LE. 0.) GO TO 80
C
C  CHECK TO MAKE SURE THAT THERE IS SLUSH AT THE
C  BEGINNING OF THE NODE-- IF NOT, JUMP TO LIQUID ROUTINES
C
      P2 = P - DP1
      IF (P2 .LE. 1.3) GO TO 198
      GO TO 199
198  WRITE(9,900)
      WRITE(9,901)
900  FORMAT(/' EXIT PRESSURE IS LESS THAN 1.1 PSIA- THE TRIPLE
      * POINT OF HYDROGEN')
901  FORMAT(/' RESET INITIAL PRESSURE AND TRY AGAIN.' )
      GO TO 70
199  CONTINUE
      EP1(J+1) = P2
      IF (P1 .GE. 190.) GO TO 70
C
C  P2 = FINAL (DOWNSTREAM) PRESSURE
C
      PAVG = (P2 + P)/2.0
C
      RHONEW = DENLPC(TOLD,P2,ID)
      RHOMIX = X2*RHOS + (1.-X2)*RHONEW
      VELFA = MDOT/(RHOMIX*AREA*3600.)
      QFLOWA = (VELFA**2-VELI**2)/(2.*GC*778.)
      QZ = (Z2-Z1)/778.
      HLIQI = ENTLPC(TOLD,P,ID)
      HFUS = 25.02
      HMIXI = -X1*HFUS + HLIQI
      HMIXF = HMIXI + QINTOT/MDOT - QZ - QFLOWA - WS
      HLIQF = ENTLPC(TOLD,P2,ID)
C
C
C  RHOMIX  MIXTURE DENSITY, LB/CU.FT
C  QFLOWA  = KINETIC ENERGY TERM, BTU/LB
C  QZ      = POTENTIAL ENERGY TERM, BTU/LB
C  HLIQI   = INITIAL ENTHALPY OF THE LIQUID AT T & P , BTU/LB
C  HFUS    = HEAT OF FUSION, BTU/LB
C  HMIXI   = INITIAL ENTHALPY OF THE MIXTURE, BTU/LB
C  HMIXF   = FINAL MIXTURE ENTHALPY, BTU/LB
C  WS      = SHAFT WORK, BTU/LB
C
C-----
C
      IF (HLIQF .LE. HMIXF) GO TO 80
C
C  CHECK TO SEE THAT THE ENTHALPY BALANCES
C
      QUALE = (HLIQF-HMIXF)/HFUS
      DQUALP = (144.*DP1A)/(778.*RHOMIX*HFUS)
      QUAL = QUALE - DQUALP
      EQUAL(J+1) = QUAL
C
C-----
C
      QUALE  FINAL SOLID FRACTION AFTER HEAT LEAKS HAVE BEEN
      TAKEN INTO ACCOUNT
      DQUALP  SLUSH LOSS DUE TO FRICTION LOSSES
      QUAL    FINAL SOLID FRACTION
C
C-----

```



IF (QUAL .LE. 0.0) GO TO 80

CHECK TO SEE THAT SLUSH EXISTS AT THE NODE EXIT; IF NOT,  
JUMP TO LIQUID ROUTINES

REPEAT CALCULATIONS IN THE PROCESS OF CONVERGENCE

RHOMIX2 = QUAL\*RHOS + (1.-QUAL)\*RHONEW  
RHOLI = DENLPC(TOLD,P,ID)  
RHOMI = QUAL\*RHOS + (1.-QUAL)\*RHOLI  
VIS2 = VISLPC(TOLD,P2,ID)  
VELF = MDOT/(RHOMIX2\*AREA\*3600.)  
QFLOW = (VELF\*\*2-VELI\*\*2)/(2.\*GC\*778.)  
REY2 = DIA\*RHOMIX2\*VELF/VIS2  
RE1(J) = REY2

IF (X2 .LE. 0.0) GO TO 71  
IF (FRICT .GE. 0.0) GO TO 71  
IF (RE .GT. 1.0E+06) GO TO 71

CALL FRICT1(QUAL,F2,REY2)

GO TO 74

CALL FRICT2(F2,REY2,ED)

CONTINUE

IF (TYPE .EQ. 'PH') GO TO 888  
IF (TYPE .EQ. 'PH ') GO TO 888  
IF (TYPE .EQ. 'PH ') GO TO 888  
IF (TYPE .EQ. 'PP') GO TO 889  
IF (TYPE .EQ. 'PP ') GO TO 889  
IF (TYPE .EQ. 'PP ') GO TO 889  
GO TO 894

WS = HP\*60./(.02356\*MDOT)  
DP2A = 0.0  
GO TO 890

WS IS IN UNITS OF BTU/LB

DP2B = DPP  
DP2A = 0.0  
WSMIN = DPP\*MDOT\*144./RHOMIX  
WSMIN IS IN UNITS OF FT-LB/HR  
HP1 = WSMIN/EFFP  
W = HP1/(MDOT\*778.)  
HP = HP1/(3600.\*550.)  
GO TO 891

CONTINUE

DP2A = (RHOMIX2\*CNLD\*VELF\*\*2\*F2)/(2.0\*144.\*GC)

CONTINUE

DP2B = RHOMIX2\*QZ\*778./144. + QFLOW\*778.\*RHOMIX2/144.  
+ EFFP\*WS\*RHOMIX2\*778./144.

CONTINUE

DP2 = DP2A + DP2B  
EDP2(J+1) = DP2  
FLOWF = (MDOT\*7.48)/(RHOMIX2\*60.)

```

      HMIXF      = HMIXI + QINTOT/MDOT - QZ - QFLOW - WS
      QUALE     = (HLIQF-HMIXF)/HFUS
C
      DQUALP    = (144.*DP2A)/(778.*RHOMIX2*HFUS)
C
      QUAL      = QUALE-DQUALP
      EQUAL(J+1) = QUAL
      ERHO(J+1) = RHOMIX2
C
      RECHECK TO SEE IF SLUSH EXISTS AT THE NODE EXIT
C
      IF (QUAL .LE. 0.0) GO TO 80
C
      CHECK TO MAKE SURE THE DENSITY AND PRESSURE CONVERGE
C
      IF (ABS(DP2-DP1) .LE. 0.0001) GO TO 10
        GO TO 20
20    DP1 = DP2
      N = N+1
      IF (N .GE. 400) GO TO 15
      X2 = QUAL
      GO TO 30
C
C
10    DELRHO = ABS(RHOMIX2 - RHOMIX)
      IF ((DELRHO) .LE. 0.0001) GO TO 15
        GO TO 16
16    RHOMIX = RHOMIX2
      N1 = N1+1
      X2 = QUAL
      IF (N1 .GE. 400) GO TO 15
      GO TO 30
C
C
      -----
C
80    CONTINUE
C
      FIRST CHECK TO SEE IF SLUSH EXISTS AT THE ENTRANCE.
C
      IF (XIN .GT. 0.0) GO TO 27
        GO TO 28
27    MDOT = EMDOT(1)
28    CONTINUE
      TIN = TOLD
      T   = TOLD
C
      CALCULATE THE SATURATION TEMPERATURE AT THE INITIAL NODE
      PRESSURE. IF THE SAT. TEMP. IS EQUAL TO THE INITIAL NODE
      TEMPERATURE ASSUME SATURATED CONDITIONS AND CALCULATE
      SATURATED PROPERTIES USING APPROPRIATE GASPLUS ROUTINES.
C
      TSAT1 = TS(J)
      TSAT2 = TSAT(P,ID)
      IF (TS(J) .GE. ET(J)) GO TO 21
        IF (ABS(TSAT2-T) .LE. .05) GO TO 21
          GO TO 82
C
      ASSUME THE PRESSURE DROP IS EQUAL TO THAT INITIALLY CALCULATED.

```

```

C      P      INITIAL PRESSURE, PSI
C      PIN     FINAL NODE PRESSURE, PSI
C
21     DP      DP1
      PIN     P    DP
      RHOLIN   DLPSAT(PIN, ID)
      HLIN     HLPSAT(P, ID)
      RHOL     DLPSAT(P, ID)
      GO TO 86
82     CONTINUE

      CALCULATE LIQUID PROPERTIES
      (IF NO SLUSH EXISTS AT ENTRANCE OR EXIT)

      X2       0.0
      QVAL     0.0
      DP       DP1
      PIN     P    DP
      TSAT3    TSAT(PIN, ID)
      HLIN     ENTHPC(TIN, P, ID)
      RHOLIN   DENSPC(TIN, P, ID)
      RHOL     DENSPC(T, PIN, ID)
      HLIQL    ENTHPC(T, PIN, ID)
      CP1      CPPC(TIN, PIN, ID)
      CP2      CPPC(T, PIN, ID)
      CPLIQ    (CP1+CP2)/2.
      MDOT     FLOW*RHOL*60./7.48
      WRITE(3, 853) P, PIN, T, HLIQL, HLIN
C 853   FORMAT(/' P = ', F5.2, 2X, ' PIN = ', F6.2, 2X, ' T = ',
      F6.2, 2X, ' HLIQL = ', F7.2, 2X, ' HLIN = ', F8.2)
      CONTINUE
86     IF (J .GT. 1) MDOT=EMDOT(1)

      P      INITIAL PRESSURE, PSI
      PIN    FINAL PRESSURE

      EP1(J+1) PIN
      VEL2    MDOT/(RHOLIN*AREA*3600.)
      VELIN   MDOT/(RHOL*AREA*3600.)
      QFLOWL  (VEL2**2 - VELIN**2)/(2.*GC*778.)
      QZ      (Z2 - Z1)/(778.)
      HSAT    HLPSAT(PIN, ID)

      INCLUDE THE HEAT CAPACITY OF THE SOLID IN THE SLUSH

      IF (XIN .LE. 0.0) GO TO 90
      GO TO 91
91     HMELT   0.0
      GO TO 92
      IF (X1 .LE. 0.0) GO TO 93
      HMELT    (X1)*HFUS
      GO TO 92
92     HMELT   0.0
      CONTINUE

      QF      HLIN * QINTOT/MDOT   QZ - QFLOWL   HMELT   WS
      HF      HF
      IF (HF .GE. HSAT) GO TO 83
      FI      TEMPPC(HF, PIN, ID)

```

```

      IF (TL .LE. 0.) GO TO 83
      GO TO 84
C
C      IF (T > TSAT) ASSUME T=TSAT AND MAKE A COMMENT
C
83      TL      = TSAT(PIN, ID)
      TS(J+1) = TL
      HF      = HSAT
      HF1(J+1) = HSAT
      RHOL    = DLPSAT(PIN, ID)
84      RHONEW  = RHOL
      VIS2    = VIS1
      DP2     = DP
      HLIQ1   = HLIN
      HLIQF   = HF
      REY2    = RE
      P1      = P
      P2      = PIN
      F2      = F
      VELI    = VELIN
      VELF    = VEL2
      QFLOW   = QFLOWL

      SET OUTPUT TEMPERATURE AND PRESSURE EQUAL TO INITIAL
      TEMP. & PRESSURE OF NEXT NODE

      ET(J+1) = TL
      EDP2(J+1) = DP
      IF (ET(J+1) .LT. 25.20) ET(J+1)=25.20
C
C      CHECK TEMPERATURE CONVERGENCE
C
      IF ((TL-T) .GE. 0.01) GO TO 81
      GO TO 15
81      N2 = N2 + 1
      IF (N2 .GE. 400) GO TO 15
      T = TL
      IF (HF .GE. HSAT) GO TO 88
      GO TO 82
88      RHOLIN = DLPSAT(PIN, ID)
      HLIQL   = HSAT
      GO TO 86
15      CONTINUE
      IF (LONG .LT. 0.0) GO TO 32
      CONTINUE
C
C      WRITE LONG, INDIVIDUAL VARIABLE OUTPUT IF LONG > 0
C
C      WRITE(8,146) TYPE
146      FORMAT(/' ELEMENT TYPE = ',A4)
C
      WRITE(8,100) QINTOT, FLOW, E, DIAM
      WRITE(8,141) OUT, LONG, XPUT, FRICT, HEAT
      WRITE(8,142) TAMB, XRS, VAC, DIN, DOUT
      WRITE(8,109) JA, CNLD, RHOS, RHONEW
      WRITE(8,101) T, P, X1, LEN
      WRITE(8,102) MDOT, RHOMIX2
      WRITE(8,106) VIS2, HLIQ1

```

```

WRITE(8,103) HMIXI,HMIXF
WRITE(8,107) REY2,F2,HLIQF
WRITE(8,115) DP1A,DP1B,DP2A,DP2B
WRITE(8,104) P2,DP2,QUAL
WRITE(8,108) N,N1,PIN1
WRITE(8,110) QUALE,DQUALP
WRITE(8,111) Z2,Z1,QZ
WRITE(8,112) VELI,VELF,QFLOW
WRITE(8,113) HF,TL,HLIQL,N2
WRITE(8,114) HSAT,TIN,HLIN
WRITE(8,145) WS,EFFP,DPP,HP
100  FORMAT(/' QIN = ',F10.3,1X,' BTU/HR',2X,' FLOW = ',F9.2,1X,
      ' GPM',1X,' E = ',F8.6,1X,' DIA = ',F7.3,1X,' IN.')
141  FORMAT(/' OUT = ',F5.2,2X,' LONG = ',F5.2,2X,' XPUT = ',
      F5.2,2X,' FRICT = ',F5.2,2X,' HEAT = ',F5.2)
142  FORMAT(/' TAMB = ',F6.2,2X,' XRS = ',F4.1,2X,' VAC = ',
      F5.2,2X,' DIN = ',F6.2,2X,' DOUT = ',F6.2)
101  FORMAT(/' T = ',F7.2,1X,' R',2X,' PI = ',F8.3,1X,' PSIA',2X,
      ' XI = ',F7.4,2X,' LENGTH = ',F10.4,1X,' FT')
108  FORMAT(/' NODE = ',I4,2X,' L/D = ',F8.3,2X,
      ' RHOS = ',F9.4,2X,' RHOL = ',F9.4,1X,' LB/CU.FT')
102  FORMAT(/' MDOT = ',F9.2,1X,' LB PER HR',2X,' RHO,MIX = ',F8.3,1X,
      ' LB PER CU.FT.')
106  FORMAT(/' MU, MIX = ',1PE14.7,1X,' LB PER FT-S',3X,
      ' HLIQI = ',OPF9.3,1X,' BTU PER LB')
103  FORMAT(/' HMIX,I = ',F9.3,1X,' BTU PER LB',3X,' HMIX,F = ',F9.3,
      ' BTU/LB')
107  FORMAT(/' RE = ',1PE14.7,2X,' F = ',OPF10.7,2X,' HLIQF = ',F9.3)
115  FORMAT(/' DP1A = ',F8.4,2X,' DP1B = ',F8.4,2X,' DP2A = ',F8.4,2X,
      ' DP2B = ',F8.4)
104  FORMAT(/' PF = ',F8.3,1X,' PSIA',2X,' DP2 = ',F9.5,1X,' PSID',2X,
      ' SOLID FRACTION = ',F10.7)
105  FORMAT(/' N = ',I3,2X,' N1 = ',I3,2X,' PIN = ',F9.5)
109  FORMAT(/' QUALE = ',F10.6,2X,' DQUALP = ',F10.6)
111  FORMAT(/' Z2 = ',F7.2,3X,' Z1 = ',F7.2,3X,' QZ = ',F9.3,1X,
      ' BTU/LB')
112  FORMAT(/' VI = ',F8.3,2X,' VF = ',F8.3,1X,' FT/S',2X,
      ' QFLOW = ',1PE14.7,2X,' BTU/LB')
113  FORMAT(/' HF = ',F9.3,2X,' TLIQ = ',F7.2,1X,' R',
      ' HLIQL = ',F9.3,2X,' N2 = ',I3)
114  FORMAT(/' HSAT = ',F9.3,2X,' TIN = ',F7.2,2X,' HLIN = ',F9.3,
      ' BTU/LB')
145  FORMAT(/' WS = ',F10.3,2X,' EFFP = ',F5.3,2X,' DPP = ',F8.3,
      ' HP = ',F8.3)
WRITE(8,105)
105  FORMAT(/'.....
      '.....')

```

PRINT OUT INDIVIDUAL NODE CONDITIONS OR END NODE CONDITIONS

```

CONTINUE
IFNEW = 0
IFP2(1) = 0
DO 26 J = 1,NUM+1
  EMDOT(J) = 0
  IF (EQUAL(J),LE,0.0) GO TO 85
  EMDOT(M+1) = EMDOT(NUM)

```

```

      MPHR = MDOT/3600.
      EPNEW = EDP2(J) + EPNEW
      DELQ = XIN - EQUAL(J)
      DELTP = PIN1 - EP1(J)
      IF (OUT .LT. 0.0) GO TO 36
      WRITE(9,200) NODE(J), EPNEW, MPHR, EQUAL(J), DELQ, EP1(J)
200  FORMAT(I3,3X,F10.5,5X,F8.4,9X,F7.4,9X,F7.4,4X,F10.4)
      IF ((RE1(J) .GT. 1.0E+06) .AND. (FRICT .LT. 0.)) GO TO 72
      GO TO 703
72  WRITE (9,701)
701  FORMAT(/' RE > 1.0E+06 - RE LIMIT EXCEEDED ')
703  CONTINUE
      WRITE(9,201)
201  FORMAT('-----')
      GO TO 87
C
C
85  CONTINUE
      EMDOT(NUM+1)=EMDOT(NUM)
      MPHR=MDOT/3600.
      EPNEW = EDP2(J) + EPNEW
      IF (OUT .LT. 0.0) GO TO 36
      WRITE(9,202) NODE(J), EPNEW, MPHR, QUAL, ET(J), EP1(J)
202  FORMAT(I3,3X,F10.5,5X,F8.4,9X,F7.4,6X,F10.3,6X,F10.4)
      IF (HF1(J) .GE. HSAT) GO TO 98
      GO TO 99
98  IF (J .LT. 2) GO TO 99
      WRITE (9,204)
204  FORMAT(' TEMPERATURE EXCEEDS SATURATION TEMPERATURE')
99  WRITE(9,203)
203  FORMAT('-----')
      GO TO 87
87  CONTINUE
36  CONTINUE
C
C
      IF (OUT .GE. 0.0) GO TO 56
      IF (QUAL .LE. 0.0) GO TO 76
      WRITE(9,600) FLOW, MPHR, QUAL, DELQ, EPNEW
600  FORMAT(1X,F10.3,5X,F8.3,9X,F7.4,11X,F7.4,7X,F10.4)
      IF ((REY2 .GT. 1.0E+06) .AND. (FRICT .LT. 0.)) GO TO 37
      GO TO 77
37  WRITE (9,702)
702  FORMAT(/' RE > 1.0E+06 - RE LIMIT EXCEEDED ')
77  CONTINUE
      WRITE(9,601)
601  FORMAT(3X,'-----')
      GO TO 56
C
C
76  CONTINUE
      IF (TL .LT. 25.2) TL = 25.2
      WRITE(9,602) FLOW, MPHR, QUAL, TL, EPNEW
602  FORMAT(1X,F10.3,5X,F8.3,9X,F7.4,8X,F10.3,8X,F10.4)
      IF (HF .GE. HSAT) GO TO 78
      GO TO 79
78  WRITE (9,604)
604  FORMAT(/10X,' TEMPERATURE EXCEEDS SATURATION TEMPERATURE')

```

```

79  WRITE(9,603)
603  FORMAT(3X,'-----
      '-----')
C
55  CONTINUE
C
C  -----
C  RESET FLOW RATE AND REPEAT CALCULATIONS
C  -----
      IF (FLOINC .LE. 0.0) GO TO 63
      IF (FLOW .LT. FLOMAX) GO TO 60
      GO TO 70
60  FLOW = FLOW + FLOINC
      GO TO 50
63  WRITE(9,999)
64  FORMAT(' FLOW INCREMENT, FLOINC, WAS NOT INCLUDED
      IN THE INPUT - REVIEW INPUT DATASET')
70  CONTINUE
      END

```

SUBROUTINE FRICT1(X,F,RE)

-----  
THIS PROGRAM CALCULATES THE FRICTION FACTOR AS A  
FUNCTION OF REYNOLDS NUMBER FOR SLUSH HYDROGEN.  
THIS DATA WAS OBTAINED USING A SMOOTH, VACUUM-JACKETED  
PIPE 0.652 INCH IN DIAMETER AS DESCRIBED IN  
SINDT, LUDTKE, AND DANNEY,  
NBS TECHNICAL NOTE 377,  
"SLUSH HYDROGEN FLUID CHARACTERIZATION AND  
INSTRUMENTATION," FEB. 1969. DATA WAS CURVE-FITTED  
TO OBTAIN THE EQUATIONS IN THIS PROGRAM.

NOTES: THE GEOMETRIC CURVE FIT, FG, TENDS TO GIVE  
THE BEST FIT OVER THE ENTIRE RANGE; THIS IS EXPECTED  
AS THE FRICTION FACTOR NORMALLY TAKES THE FORM  
 $F = A/RE^{.25}$

STANDARD FRICTION FACTORS GIVEN IN THE LITERATURE

ARE, FOR TURBULENT FLOW,

$F = 4 \cdot 0.079 \cdot RE^{-.25}$  AND

$F = 4 \cdot 0.046 \cdot RE^{-.20}$

THE SECOND EQUATION GIVES A BETTER FIT TO EXPERIMENTAL  
DATA. IN ADDITION, THE EQUATION OFTEN USED IS

$F \cdot 0.5 = 4.06 \cdot \log(RE/F \cdot 0.5) - 0.60$

THIS IS USED TO COMPARE DATA IN THE NBS DATA.

FOR LAMINAR FLOW,  $F = 64/RE$

-----  
THE FRICTION FACTOR CALCULATIONS CAN ONLY BE USED FOR  
 $RE \geq 1E06$

1 GEOMETRIC FIT ( $F = AX^{.B}$ ) OF THE FRICTION FACTOR

```

      IF (X .GE. 0.5 ) GO TO 1
        GO TO 2
2     IF (X .GE. 0.4) GO TO 3
        GO TO 4
4     IF (X .GE. 0.3) GO TO 5
        GO TO 6
6     IF (X .GE. 0.2) GO TO 7
        GO TO 7
1     FG = 3.0775*RE**-.41267
      F = FG
      GO TO 10
3     FG4 = 1.94175*RE**-.38524
      FG5 = 3.0775*RE**-.41267
      F = FG4*(.5-X)*10. + FG5*(1.-(.5-X)*10.)
      GO TO 10
5     FG4 = 1.94175*RE**-.38524
      FG3 = 0.70205*RE**-.309898
      F = FG3*(.4-X)*10. + FG4*(1.-(.4-X)*10.)
      GO TO 10
7     FG2 = 0.421259*RE**-.27205
      FG3 = 0.70205*RE**-.309898
      IF (X .LE. 0.2) GO TO 11
        GO TO 13
11    AX = .2
      GO TO 12
13    AX = X
12    F = FG2*(.3-AX)*10. + FG3*(1.-(.3-AX)*10.)
10    CONTINUE
      RETURN
      END

```

C  
C

```

SUBROUTINE VACUUM (QRAD,TAMB,XRS,DIN,DOUT,TIN,E1,E2)
  REAL L,XMFP,MU

```

C

```

C-----
C  THIS PROGRAM CALCULATES THE HEAT TRANSFER THROUGH
C  VACUUM JACKETED LINES AS A FUNCTION OF OUTSIDE
C  TEMPERATURE AND PIPE DIAMETER FOR SLUSH HYDROGEN SYSTEMS.
C-----

```

C VARIABLES:

```

C      E1----EMISSIVITY OF INSIDE SURFACE,UNITLESS
C      E2----EMISSIVITY OF OUTER SURFACE,UNITLESS
C      DIN---OUTER DIAMETER OF THE INSIDE PIPE, IN
C      DOUT--INNER DIAMETER OF THE OUTER PIPE, IN
C      TAMB--AMBIENT TEMPERATURE, R
C      TIN---FLUID TEMPERATURE, R
C      XRS---NUMBER OF RADIATION SHIELDS
C      QRAD--RADIATION HEAT TRANSFER, BTU/FT-HR
C      F12---CONFIGURATION FACTOR
C      FE----EMISSIVITY FACTOR
C      SIG---STEFAN-BOLTZMANN CONSTANT
C-----

```

C CALCULATE HEAT LOSS DUE TO RADIATION

```

C      T2 = TAMB
C      T1 = TIN

```

C

10 CONTINUE



```

      IF (XRS .GT. 0.) GO TO 15
      GO TO 16
15    F12 = 1.0
      E1IN = 1./E1
      E2IN = 1./E2
      ES = 0.03
      ES = ALUMINUM SHIELD EMISSIVITY
      ESIN = 1./ES
      CON1 = E1IN + ESIN - 1.
      CON2 = (XRS-1.)*(2.0*ESIN - 1.)
      CON3 = E2IN + ESIN - 1.
      FEIN = CON1 + CON2 + CON3
      FECALC = 1./FEIN
      FEACT = 5.0*FECALC
      FE = FEACT
      GO TO 17
16    F12 = 1.0
      A1A2 = DIN/DOUT
      E1IN = 1./E1
      E2IN = 1./E2
      FEIN = E1IN + A1A2*(E2IN-1.0)
      FE = 1.0/FEIN
17    P1 = 3.14159
      SIG = 0.173E-08
      A1 = P1*DIN/12.
      TP2 = T2**4.
      TP1 = T1**4.
      QRAD = F12*FE*SIG*A1*(TP2 - TP1)
      QRAD HAS UNITS OF BTU/HR-FT
20    RETURN
      END

SUBROUTINE FRICT2(F2,RE,ED)

THIS SUBROUTINE CALCULATES FRICTION FACTORS FOR
TURBULENT FLOW USING THE COLBROOK EQUATION.

      WRITE(14,201) RE,ED,F2
201    FORMAT(1PE15.8,2X,E15.8,2X,E15.8)
      F1 = 1.325/(ABS(ALOG(ED/3.7 + 5.74/RE**0.9))**2.0)
      N = 0
      F2A = 0.021
      CONTINUE
      CON1 = (ED/3.7 + 2.51/(RE*SQRT(F2A)))
      CONF2 = 2.*ALOG10(CON1)
      F2B = 1.0/(CONF2)**2.
      IF (ABS(F2B-F2A) .LE. 0.000001) GO TO 20
      F2A = F2B
      N = N+1
      WRITE(15,203) RE,N
203    FORMAT(1PE15.8,2X,I5)
      IF (N .GT. 400) GO TO 23
      GO TO 30
20    CONTINUE
      WRITE(9,206) RE,F2,CON1
206    FORMAT(1PE15.8,2X,E15.8,2X,E15.8)

```

```

      F2 = F2B
23      RETURN
      END

C
      SUBROUTINE QFLUX(QFLA,QLA)
C
C      PROGRAM TO CALCULATE THE HEAT TRANSFER RATE
C      TO SLUSH HYDROGEN BEING TRANSFERRED IN
C      PIPING WITHOUT VACUUM-JACKETING.
C
C      QFLUX ALLOWS UP TO 3 LAYERS OF INSULATION AROUND
C      A BARE PIPE
C
C      TERRY HARDY      12-16-88
C
      DATA ID/'PH2'/
      DIMENSION NELEM(90),
      . TK(10,90),
      . INSUL(90),HIN(90),HOUT(90),U(90),QFLX(90),AREA(90),
      . VELI(90),QL(90)

      COMMON /FLUX/ EPOD(90),ETI1(90),ETI2(90),ETI3(90),
      . NNOVJ(90),NOJAC,NXMAT(20,90),
      . EXLEN(90),EXDIA(90),EXT(90),TAMB,P,X1,FLOV,EPID(90),K
C
C
      CHARACTER*10 TIM,TONE,MIM,MONE
      INTEGER INSUL,NXMAT
      REAL MDOT

C
      I = 1
      NUM = 1
      AIRV = 0.01

C
      CALL DATE(TIM)
      CALL TIME(TONE)
      WRITE(10,130) TIM,TONE
C 130  FORMAT('/' , A10,2X,A10)
C
C
      DO 54 I = 1,NUM
      IF (NXMAT(2,K) .LE. 0.0) GO TO 71
      IF (NXMAT(3,K) .LE. 0.0) GO TO 72
      IF (NXMAT(4,K) .LE. 0.0) GO TO 73
      GO TO 74
71      INSUL(I) = 1
      GO TO 75
72      INSUL(I) = 2
      GO TO 75
73      INSUL(I) = 3
      GO TO 75
74      INSUL(I) = 4
75      CONTINUE
54      CONTINUE

C
C
C      INITIALIZE PROPERTIES
C

```

```

DO 55 I = 1,NUM
TFL = EXT(K)
RHOL = DENLPC(TFL,P,ID)
RHOS = DNSOPC(TFL,P,ID)
RHOMIX = X1*RHOS + (1.-X1)*RHOL
PI = 3.14159
AREA(I) = PI*EPID(K)**2/4.0
MDOT = FLOV*RHOMIX*60./7.48
VELI(I) = MDOT/(AREA(I)*RHOMIX*3600.)
CONTINUE
DO 50 I = 1,NUM
DO 51 J = 1,INSUL(I)

```

.....  
MATCH MATERIAL CODE TO THERMAL CONDUCTIVITY VALUES  
.....

```

IF (NXMAT(J,K) .LE. 101) GO TO 11
IF (NXMAT(J,K) .EQ. 102) GO TO 12
IF (NXMAT(J,K) .EQ. 103) GO TO 13
IF (NXMAT(J,K) .EQ. 104) GO TO 14
IF (NXMAT(J,K) .EQ. 201) GO TO 21
IF (NXMAT(J,K) .EQ. 202) GO TO 22
IF (NXMAT(J,K) .EQ. 203) GO TO 23
IF (NXMAT(J,K) .EQ. 204) GO TO 24
IF (NXMAT(J,K) .EQ. 205) GO TO 25
IF (NXMAT(J,K) .EQ. 301) GO TO 31
IF (NXMAT(J,K) .EQ. 302) GO TO 32
GO TO 11

```

```

11 TK(J,K) = 0.112
GO TO 70
12 TK(J,K) = 2.717
GO TO 70
13 TK(J,K) = 22.31
GO TO 70
14 TK(J,K) = 156.0
GO TO 70
15 TK(J,K) = 0.020
GO TO 70
16 TK(J,K) = 0.032
GO TO 70
17 TK(J,K) = 0.019
GO TO 70
18 TK(J,K) = 0.019
GO TO 70
19 TK(J,K) = 0.021
GO TO 70
20 TK(J,K) = 0.014
GO TO 70
21 TK(J,K) = 0.015
GO TO 70
22 CONTINUE
23 CONTINUE
24 CONTINUE

```

THE FOLLOWING MATERIAL CODES APPLY

CODE	MATERIAL
------	----------

```

C
C      101          TEFLON
C      102          304 STAINLESS STEEL
C      103          CARBON STEEL
C      104          6063 ALUMINUM
C      201          FOAM GLASS
C      202          SILICA
C      203          POLYURETHANE
C      204          POLYSTYRENE
C      205          FOAM RUBBER
C      301          FIBER GLASS
C      302          PERLITE
C
C      THE VALUES OF THERMAL CONDUCTIVITY WERE OBTAINED
C      FROM "CRYOGENIC SYSTEMS", 2ND ED., RANDALL F. BARRON,
C      OXFORD UNIVERSITY PRESS, 1985.

```

```

C      -----
C      CALCULATE CONDUCTION HEAT TRANSFER PARAMETERS
C      -----

```

```

C      DO 60 I = 1, NUM
C          DR1 = (EPOD(K) - EPID(K))/2.0
C          COND1 = (DR1/TK(1,K))/((2.0*PI*EXLEN(K)*DR1)/
C              (ALOG(EPOD(K)/EPID(K))))
C      IF (INSUL(I) .LE. 1) GO TO 90
C
C          R1 = EPOD(K)/2.0
C          R2 = EPOD(K)/2.0 + ETI1(K)
C          R3 = R2 + ETI2(K)
C          R4 = R3 + ETI3(K)
C
C      81      DR2 = R2 - R1
C          COND2 = (DR2/TK(2,K))/((2.0*PI*EXLEN(K)*DR2)/
C              (ALOG(R2/R1)))
C          IF (INSUL(I) .LE. 2) GO TO 80
C      82      DR3 = R3 - R2
C          COND3 = (DR3/TK(3,K))/((2.0*PI*EXLEN(K)*DR3)/
C              (ALOG(R3/R2)))
C          IF (INSUL(I) .LE. 3) GO TO 80
C      83      DR4 = R4 - R3
C          COND4 = (DR4/TK(4,K))/((2.0*PI*EXLEN(K)*DR4)/
C              (ALOG(R4/R3)))
C
C      80      CONTINUE
C          CONINS = COND2 + COND3 + COND4
C      90      CONTINUE
C          CONTOT = COND1 + CONINS
C          PAMB = 14.7
C          TFL = EXT(K)
C
C      CALL SUBROUTINES FOR CALCULATION OF CONVECTIVE
C      HEAT TRANSFER COEFFICIENTS
C
C      CALL INHCO(X1,TFL,P, VELI(I),EXLEN(K),EPID(K),HIN(I))
C      CALL OUTHCO(TAMB,PAMB,AIRV,EXLEN(K),EPOD(K),HOUT(I))

```

```

HEATI = 1.0/(PI*EPID(K)*EXLEN(K)*HIN(I))
HEATO = 1.0/(PI*EPOD(K)*EXLEN(K)*HOUT(I))
UINV = HEATI + CONTOT + HEATO
U(I) = 1.0/UINV
QFLX(I) = U(I)*(TAMB-TFL)
QL(I) = QFLX(I)/EXLEN(K)
QFLA = QFLX(I)
QLA = QL(I)

```

```

-----
QFLX = HEAT TRANSFER, BTU/HR
QL = HEAT TRANSFER, BTU/HR-FT
HIN = INSIDE HEAT TRANSFER COEFFICIENT,BTU/HR-FT**2-R
HOUT = OUTSIDE HEAT TRANSFER COEFFICIENT,BTU/HR-FT**2-R
U = OVERALL HEAT TRANSFER COEFFICIENT
COND1= CONDUCTION THRU PIPE = (RO-RI)/K-A,LM
COND2= CONDUCTION THRU INSULATION MATERIAL 1
COND3= CONDUCTION THRU INSULATION MATERIAL 2
COND4= CONDUCTION THRU INSULATION MATERIAL 3
-----

```

# WRITE OUTPUT FOR INDIVIDUAL VARIABLES

```

WRITE(10,910)
FORMAT(/'
-----')
WRITE(10,911) K, EPID(K), EPOD(K), EXLEN(K)
WRITE(10,920) FLOW,AIRV,TAMB,TFL
WRITE(10,921) RHOL,RHOS,MDOT,VELI(I)
WRITE(10,912) INSUL(I), NXMAT(1,K),NXMAT(2,K),NXMAT(3,K),
. NXMAT(4,K)
WRITE(10,913) TK(1,K),TK(2,K),TK(3,K),TK(4,K)
WRITE(10,914) R1,R2,R3,R4
WRITE(10,915) COND1,COND2,COND3,COND4
WRITE(10,916) HIN(I),HOUT(I),U(I),QFLX(I)
911 FORMAT(/' ELEM = ',I5,2X,' ID = ',F8.4,2X,' OD = ',F8.4,
. 2X,' LENGTH = ',F8.3)
920 FORMAT(/' FLOW = ',F8.3,2X,' VAIR = ',F8.4,2X,' TAMB = ',
. F7.2,2X,' TFLUID = ',F7.2)
921 FORMAT(/' RHOL = ',F7.4,2X,' RHOS = ',F7.4,2X,' MDOT = ',
. F8.3,2X,' VFLUID = ',F8.4)
912 FORMAT(/' NINS = ',I4,2X,' PIPE = ',I5,2X,' MAT2 = ',I5,2X,
. ' MAT3 = ',I5,2X,' MAT4 = ',I5)
913 FORMAT(/' K1 = ',F9.4,2X,' K2 = ',F8.5,2X,' K3 = ',
. F8.5,2X,' K4 = ',F8.5)
914 FORMAT(/' R1 = ',F8.4,2X,' R2 = ',F8.4,2X,' R3 = ',F8.4,2X,
. ' R4 = ',F8.4)
915 FORMAT(/' CON1 = ',1PE14.7,2X,' CON2 = ',OPF8.5,2X,
. ' CON3 = ',F8.4,2X,' CON4 = ',F8.4)
916 FORMAT(/' HI = ',F8.4,2X,' HO = ',F8.4,2X,' U = ',F8.5,
. 2X,' QFLX = ',F12.4)

```

# CALCULATE HEAT TRANSFER FOR A NEW ELEMENT

```

CONINS = 0.0
COND1 = 0.0

```

```

COND2 = 0.0
COND3 = 0.0
COND4 = 0.0
C
60  CONTINUE
    RETURN
    END
C
C
SUBROUTINE INHCO(X,T,P,V,XLEN,DIA,HIN)
C
C  SUBROUTINE TO CALCULATE INSIDE HEAT TRANSFER COEFFICIENTS
C
CALL HYDRO(X,T,P,CP,TK,RHO,VIS)
C
C  X      SOLID FRACTION
C  T      FLUID TEMPERATURE, R
C  P      AVERAGE FLUID PRESSURE, PSI
C  V      FLUID VELOCITY, FT/SEC
C  XLEN=  FLOW LENGTH, FT
C  DIA =  FLOW DIAMETER, FT
C  CP =   SPECIFIC HEAT, BTU/LB-R
C  TK =   THERMAL CONDUCTIVITY, BTU/FT-SEC-R
C  RHO =  SLUSH DENSITY, LB/CU.FT.
C  VIS =  VISCOSITY, LB/FT-SEC
C  RE =   REYNOLD'S NUMBER
C  PR =   PRANDTL NUMBER
C  HIN =  INSIDE HEAT TRANSFER COEFFICIENT, BTU/HR-SQ.FT.-R
C
C  RE = DIA*RHO*V/VIS
C  PR = CP*VIS/TK
C  IF (RE .LT. 2100.) GO TO 10
    GO TO 20
10  HINS = 1.86*(TK/XLEN)*(RE*PR*DIA/XLEN)**0.333
    GO TO 30
20  HINS = 0.023*(TK/XLEN)*(RE**0.8)*(PR**0.333)
30  HIN = HINS*3600.
    RETURN
    END
C
C
SUBROUTINE OUTHCO(T,P,V,XLEN,DIA,HOUT)
C
C  SUBROUTINE TO CALCULATE OUTSIDE HEAT TRANSFER COEFFICIENTS
C
CALL AIR(T,P,CP,TK,RHO,VIS,BETA)
C
C  T      AIR TEMPERATURE, R
C  P      AVERAGE AIR PRESSURE, PSI
C  V      AIR VELOCITY, FT/SEC
C  XLEN=  FLOW LENGTH, FT
C  DIA =  FLOW DIAMETER, FT
C  CP =   SPECIFIC HEAT, BTU/LB R
C  TK =   THERMAL CONDUCTIVITY, BTU/FT SEC R
C  RHO =  AIR DENSITY, LB/CU.FT.
C  VIS =  AIR VISCOSITY, LB/FT SEC
C  RE =   REYNOLDS NUMBER
C  PR =   PRANDTL NUMBER
C  GR =   GRASHOF NUMBER
C  HOUT=  OUTSIDE HEAT TRANSFER COEFFICIENT, BTU/HR-SQ.FT.-R

```

```

C
G - 32.2
RE = DIA*RHOMIX*V/VIS
PR = CP*VIS/TK
GR = XLEN**3*RHO**2*G*BETA*(T)/(VIS**2)
C
NOTE: THE APPROXIMATION FOR BETA IS 1/TAVG = 1/T
C
IF (V .LT. 0.1) GO TO 15
IF (RE .LT. 2100.) GO TO 10
GO TO 20
15 CB = 0.333
CA = 0.140
HOUTS = (TK/XLEN)*CA*(GR*PR)**CB
GO TO 30
10 HOUTS = 1.86*(TK/XLEN)*(RE*PR*DIA/XLEN)**0.333
GO TO 30
20 HOUTS = 0.023*(TK/XLEN)*(RE**0.8)*(PR**0.333)
30 HOUT = HOUTS*3600.
RETURN
END

```

```

SUBROUTINE AIR(T,P,CP,TK,RHO,VIS,BETA)

```

```

DATA ON AIR OBTAINED FROM GASPLUS

```

```

DATA ID/'AIR'/
CP = CPVPC(T,P,ID)
TK = TCNVPC(T,P,ID)
RHO = DENVPC(T,P,ID)
VIS = VISVPC(T,P,ID)
BETA = 1.0/T

```

```

RETURN
END

```

```

SUBROUTINE HYDRO(X,T,P,CP,TK,RHO,VIS)

```

```

DATA ON HYDROGEN OBTAINED FROM GASPLUS

```

```

DATA ID/'PH2'/
CP = CPLPC(T,P,ID)
TK = TCNLPC(T,P,ID)
RHOL = DENLPC(T,P,ID)
VIS = VISLPC(T,P,ID)
RHOS = DNSOPC(T,P,ID)
RHO = X*RHOS + (1. - X)*RHOL

```

```

RETURN
END

```

## APPENDIX C

### NOTES FOR NASA LEWIS VAX USERS

The following steps are required to use the FLUSH code on the NASA Lewis VAX system:

(1) Create a file called FLUSH.INP as described in the users guide. Several points could be of help in producing this file:

- Namelists must always start in column 2 of your file. Therefore, the statements \$PARAMS and \$SEND must start in column 2.
- The element variable ELE in the element description section is I4 format. Therefore, single-digit integers need to be in column 4; two-digit variables need to start in column 3. TYPE is A4 format and can be placed in columns 5 to 8. Because the remaining variables (DIA, LENGTH, TEMP Q, K, and HT) have F10.2 format, they need to be placed in lines 9 to 18, 19 to 28, etc.
- The line with the descriptors ELE, DIA, etc., is required for the code to work, but because this is A80 format, the user can vary the actual names of the descriptors, if desired.
- If the radiation heat transfer parameters are to be defined, a blank line is required between the element description section and the radiation heat transfer parameter section.

(2) To run the code, issue the statement

```
RUN [SPTELH.FLUSH]FLUSH
```

(3) Output will be set to unit 9, in a file called FOR009.DAT, for the final output conditions or the individual node output. The longer, individual variable output will be sent to unit 8 if LONG is greater than 0.

(4) Note that to run FLUSH the user needs access to the GASPLUS properties code. To do this, simply add the following line to the file LOGIN.COM;

```
ASSIGN NASA$SPID:[SPANK.GASPLUS]GASPLUS.OLB LNK$LIBRARY
```



## REFERENCES

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3. Ewart, R.O.; and Dergance, R.H.: Cryogenic Propellant Densification Study. (MCR-78-586, Martin Marietta Corp.; NASA Contract NAS3-21014) NASA CR-159438, 1978.
4. Fowler, J.R.: GASPLUS User's Manual. NASP CR-1012, Mar. 1988.
5. Barron, R.F.: *Cryogenic Systems*. 2nd ed., Oxford University Press, Oxford, 1985.
6. Ludtke, P.R.; and Voth, R.O.: A Study of LC-39 Cryogenic Systems. (NBS Report 10705, National Bureau of Standards, NASA CR-126386, 1971.
7. *Flow of Fluids Through Valves, Fittings, and Pipe*. Crane Co., Hydro-Aire Division, Chicago, IL, TP-410, 1965.
8. Siegel, R.; and Howell, J.R.: *Thermal Radiation Heat Transfer*. 2nd ed., McGraw-Hill, New York, 1981.
9. Hannum, N.P.: Technology Issues Associated With Fueling The National Aerospace Plane With Slush Hydrogen. NASA TM-101386, 1988.
10. Sindt, C.: A Summary of the Characterization Study of Slush Hydrogen. *Cryogenics*, vol. 10, no. 5, Oct., 1970, pp. 372-380.
11. Voth, R.O., et al: On Process Design Criteria, Data and Methodology Related to Slush Hydrogen Technology. NASP TM-1006, 1987.

TABLE 1. - ELEMENT TYPE FOR USE IN  
INPUT FILE

Type	Corresponding element
PI	Straight pipe
BE	Bellows
BA	Bayonet fitting
FX	Flexible tubing
EL	90° Elbow
E4	45° Elbow
YV	Y-pattern globe valve
VA	Standard globe valve
AV	Angle valve
GV	Gate valve
BV	Ball valve
PP	Pump, $\Delta P$ specified
PH	Pump, horsepower specified

TABLE 2. - EXAMPLE FLUSH CODE INPUT FILE

[Analysis of a 100-ft-long, 1.5-in.-diameter Schedule 5S Line.]

## \$PARAMS

FLOMIN=10.0, FLOMAX=100.0, FLOINC=5.0, E = 0.00015,  
 P = 35., X1 = 0.5,  
 ZIN=25., NUM=18, OUT=1.,  
 LONG=-1., XPUT=1., FRICT=+1.0, HEAT=+1.0,  
 TAMB=540., VAC=0.0, XRS=10., NOJAC=0.

## \$END

ELE	TYPE	DIA, in.	LENGTH, ft	TEMP., R.	Q	K=fL/D	HT., ft
1	PI	1.770	1.000	25.20	0.633	0.0	25.0
2	VA	1.770	0.500	25.20	27.30	7.58	25.0
3	PI	1.770	19.00	25.20	0.633	0.0	25.0
4	BE	1.770	1.000	25.20	1.899	0.0	25.0
5	BA	1.770	0.790	25.20	14.78	0.0	25.0
6	PI	1.770	19.00	25.20	0.633	0.0	25.0
7	BE	1.770	1.000	25.20	1.899	0.0	25.0
8	BA	1.770	0.790	25.20	14.78	0.0	25.0
9	PI	1.770	19.00	25.20	0.633	0.0	25.0
10	BE	1.770	1.000	25.20	1.899	0.0	25.0
11	BA	1.770	0.790	25.20	14.78	0.0	25.0
12	PI	1.770	19.00	25.20	0.633	0.0	25.0
13	BE	1.770	1.000	25.20	1.899	0.0	25.0
14	BA	1.770	0.790	25.20	14.78	0.0	25.0
15	PI	1.770	19.00	25.20	0.633	0.0	25.0
16	BE	1.770	1.000	25.20	1.899	0.0	25.0
17	VA	1.770	0.500	25.20	27.30	7.58	25.0
18	PI	1.770	1.000	25.20	0.633	0.0	25.0

TABLE 3. - HEAT LEAK ESTIMATES  
FOR PIPES CARRYING LIQUID  
HYDROGEN

[Estimates from CVI, Inc.]

Pipe diameter, in.	Rigid pipe	Flexible pipe heat leak, Btu/hr ft
0.75	0.462	1.293
1	.515	1.545
1.5	.633	1.899
2	.854	2.562
3	1.080	3.240
4	1.407	4.221
6	1.835	5.505

TABLE 4. - HEAT LEAK AND FLOW COEFFICIENT ESTIMATES  
FOR LIQUID HYDROGEN VALVES AND FITTINGS

[Estimates from Cryogenic Energy Co.]

Pipe diameter, in.	Valve heat leak, Btu/hr ft	Flow coefficient, $C_v$		Bayonet fitting heat leak, Btu/hr ft
		Y	Globe	
0.50	8.90	9.0	6.6	7.27
.75	14.10	12.0	8.0	7.27
1	21.80	22.0	14.0	10.02
1.5	27.30	47.0	34.0	14.78
2	27.30	96.0	50.0	23.19

TABLE 5. - EXAMPLE FLUSH CODE SHORT OUTPUT FILE

[Node conditions not listed.]

## Input Parameters

Element	Type	Dia., in.	Length, ft	Temp., R	Q, Btu/hr-ft	K=f L/D
1	PI	1.770	1.000	25.20	0.633	0.000
2	VA	1.770	0.500	25.20	27.300	7.580
3	PI	1.770	19.000	25.20	0.633	0.000
4	BE	1.770	1.000	25.20	1.899	0.000
5	BA	1.770	0.790	25.20	14.780	0.000
6	PI	1.770	19.000	25.20	0.633	0.000
7	BE	1.770	1.000	25.20	1.899	0.000
8	BA	1.770	0.790	25.20	14.780	0.000
9	PI	1.770	19.000	25.20	0.633	0.000
10	BE	1.770	1.000	25.20	1.899	0.000
11	BA	1.770	0.790	25.20	14.780	0.000
12	PI	1.770	19.000	25.20	0.633	0.000
13	BE	1.770	1.000	25.20	1.899	0.000
14	BA	1.770	0.790	25.20	14.780	0.000
15	PI	1.770	19.000	25.20	0.633	0.000
16	BE	1.770	1.000	25.20	1.899	0.000
17	VA	1.770	0.500	25.20	27.300	7.580
18	PI	1.770	1.000	25.20	0.633	0.000

TABLE 5. - Concluded.

## FLUSH Code Output

Pipe Diameter = 1.77 in.      Flow = 10.00 gpm  
 Total Pipe Length = 106.16 ft      Initial Solid Fraction = 0.500  
 Pipe Roughness = 0.000150      Initial Temperature = 25.20 R

Flow, gpm	Flow, lb/s	Final Solid Fraction	Change in x	Pressure Drop, psi
10.000	0.114	0.4858	0.0142	0.0300
15.000	0.171	0.4904	0.0096	0.0659
20.000	0.227	0.4926	0.0074	0.1155
25.000	0.284	0.4938	0.0062	0.1788
30.000	0.341	0.4946	0.0054	0.2558
35.000	0.398	0.4950	0.0050	0.3465
40.000	0.455	0.4952	0.0048	0.4509
45.000	0.512	0.4953	0.0047	0.5690
50.000	0.569	0.4952	0.0048	0.7007
55.000	0.626	0.4951	0.0049	0.8462
60.000	0.682	0.4949	0.0051	1.0053
65.000	0.739	0.4946	0.0054	1.1782
70.000	0.796	0.4942	0.0058	1.3647
75.000	0.853	0.4938	0.0062	1.5649
80.000	0.910	0.4933	0.0067	1.7789
85.000	0.967	0.4928	0.0072	2.0065
90.000	1.024	0.4923	0.0077	2.2479
95.000	1.081	0.4916	0.0084	2.5029
100.000	1.137	0.4910	0.0090	2.7717

TABLE 6. - EXAMPLE FLUSH CODE OUTPUT FOR NODE CONDITIONS LISTED

Pipe Diameter = 1.77 in.		Flow = 10.00 gpm			
Total Pipe Length = 106.16 ft		Initial Solid Fraction = 0.500			
Pipe Roughness = 0.000150		Initial Temperature = 25.20 R			
.....					
Node	Pressure Drop, psi	Flow, lb/s	Final Solid Fraction	Change in x	Pressure, psi
1	0.00000	0.1137	0.5000	0.0000	35.0000
2	0.00015	0.1137	0.4999	0.0001	34.9999
3	0.00725	0.1137	0.4986	0.0014	34.9928
4	0.01010	0.1137	0.4974	0.0026	34.9899
5	0.01025	0.1137	0.4972	0.0028	34.9898
6	0.01037	0.1137	0.4961	0.0039	34.9896
7	0.01323	0.1137	0.4949	0.0051	34.9868
8	0.01338	0.1137	0.4947	0.0053	34.9866
9	0.01350	0.1137	0.4936	0.0064	34.9865
10	0.01635	0.1137	0.4924	0.0076	34.9837
11	0.01650	0.1137	0.4922	0.0078	34.9835
12	0.01662	0.1137	0.4911	0.0089	34.9834
13	0.01948	0.1137	0.4899	0.0101	34.9805
14	0.01963	0.1137	0.4897	0.0103	34.9804
15	0.01975	0.1137	0.4886	0.0114	34.9803
16	0.02261	0.1137	0.4874	0.0126	34.9774
17	0.02276	0.1137	0.4872	0.0128	34.9773
18	0.02986	0.1137	0.4858	0.0142	34.9701
19	0.03001	0.1137	0.4858	0.0142	34.9700
.....					

TABLE 6. - Concluded.

## FLUSH Code Output

Pipe Diameter = 1.77 in.      Flow = 15.00 gpm  
 Total Pipe Length = 106.16 ft      Initial Solid Fraction = 0.500  
 Pipe Roughness = 0.000150      Initial Temperature = 25.20 R

Node	Pressure Drop, psi	Flow, lb/s	Final Solid Fraction	Change in x	Pressure, psi
1	0.00000	0.1706	0.5000	0.0000	35.0000
2	0.00032	0.1706	0.5000	0.0000	34.9997
3	0.01629	0.1706	0.4990	0.0010	34.9837
4	0.02242	0.1706	0.4982	0.0018	34.9776
5	0.02274	0.1706	0.4981	0.0019	34.9773
6	0.02300	0.1706	0.4973	0.0027	34.9770
7	0.02913	0.1706	0.4965	0.0035	34.9709
8	0.02945	0.1706	0.4964	0.0036	34.9705
9	0.02970	0.1706	0.4957	0.0043	34.9703
10	0.03584	0.1706	0.4949	0.0051	34.9642
11	0.03616	0.1706	0.4947	0.0053	34.9638
12	0.03641	0.1706	0.4940	0.0060	34.9636
13	0.04255	0.1706	0.4932	0.0068	34.9574
14	0.04287	0.1706	0.4930	0.0070	34.9571
15	0.04312	0.1706	0.4923	0.0077	34.9569
16	0.04926	0.1706	0.4915	0.0085	34.9507
17	0.04958	0.1706	0.4914	0.0086	34.9504
18	0.06556	0.1706	0.4904	0.0096	34.9344
19	0.06589	0.1706	0.4904	0.0096	34.9341



TABLE 7. - EXAMPLE FLUSH CODE INPUT FILE FOR HEAT LEAK CALCULATIONS ON  
ELEMENTS 1 AND 9

\$PARAMS

FLOMIN=10.0, FLOMAX=100.0, FLOINC=5.0, E = 0.00015,  
P = 35., X1 = 0.5,  
ZIN=25., NUM=18, OUT=-1.,  
LONG=-1., XPUT=1., FRICT=+1.0, HEAT=+1.0,  
TAMB=540., VAC=2.0, XRS=10., NOJAC=0.

\$END

ELE	TYPE	DIA, in.	LENGTH, ft	TEMP., R.	Q	K=fL/D	HT., ft
1	PI	1.770	1.000	25.20	-1.00	0.0	25.0
2	VA	1.770	0.500	25.20	27.30	7.58	25.0
3	PI	1.770	19.00	25.20	0.633	0.0	25.0
4	BE	1.770	1.000	25.20	1.899	0.0	25.0
5	BA	1.770	0.790	25.20	14.78	0.0	25.0
6	PI	1.770	19.00	25.20	0.633	0.0	25.0
7	BE	1.770	1.000	25.20	1.899	0.0	25.0
8	BA	1.770	0.790	25.20	14.78	0.0	25.0
9	PI	1.770	19.00	25.20	-1.00	0.0	25.0
10	BE	1.770	1.000	25.20	1.899	0.0	25.0
11	BA	1.770	0.790	25.20	14.78	0.0	25.0
12	PI	1.770	19.00	25.20	0.633	0.0	25.0
13	BE	1.770	1.000	25.20	1.899	0.0	25.0
14	BA	1.770	0.790	25.20	14.78	0.0	25.0
15	PI	1.770	19.00	25.20	0.633	0.0	25.0
16	BE	1.770	1.000	25.20	1.899	0.0	25.0
17	VA	1.770	0.500	25.20	27.30	7.58	25.0
18	PI	1.770	1.000	25.20	0.633	0.0	25.0

Radiation Heat Transfer Parameters

ELE	O.D.	I.D.	Ei	Eo
1	1.900	3.334	0.30	0.30
9	1.900	3.334	0.30	0.30

TABLE 8. - SCHEDULE 5S STAINLESS  
STEEL PIPE DATA

Nominal size, in.	Inner diameter, in.	Outer diameter, in.
0.50	0.710	0.840
.75	.920	1.050
1	1.185	1.315
1.5	1.770	1.900
2	2.245	2.375
2.5	2.709	2.875
3	3.334	3.500
4	4.334	4.500
5	5.345	5.563
6	6.407	6.625
8	8.407	8.625
10	10.482	10.750
12	12.438	12.750

TABLE 9. - EXAMPLE FLUSH CODE OUTPUT FOR HEAT LEAK CALCULATIONS  
ON ELEMENTS 1 AND 9

.....

Input Parameters

Element	Type	Dia., in.	Length, ft	Temp., R	Q, Btu/hr-ft	K f L/D
1	PI	1.770	1.000	25.20	0.552	0.000
2	VA	1.770	0.500	25.20	27.300	7.580
3	PI	1.770	19.000	25.20	0.633	0.000
4	BE	1.770	1.000	25.20	1.899	0.000
5	BA	1.770	0.790	25.20	14.780	0.000
6	PI	1.770	19.000	25.20	0.633	0.000
7	BE	1.770	1.000	25.20	1.899	0.000
8	BA	1.770	0.790	25.20	14.780	0.000
9	PI	1.770	19.000	25.20	0.552	0.000
10	BE	1.770	1.000	25.20	1.899	0.000
11	BA	1.770	0.790	25.20	14.780	0.000
12	PI	1.770	19.000	25.20	0.633	0.000
13	BE	1.770	1.000	25.20	1.899	0.000
14	BA	1.770	0.790	25.20	14.780	0.000
15	PI	1.770	19.000	25.20	0.633	0.000
16	BE	1.770	1.000	25.20	1.899	0.000
17	VA	1.770	0.500	25.20	27.300	7.580
18	PI	1.770	1.000	25.20	0.633	0.000

Radiation Heat Transfer Parameters

Element	Inner O.D.	Outer I.D.	Ei	Eo
1	1.900	3.334	0.30000	0.30000
9	1.900	3.334	0.30000	0.30000

TABLE 9. - Concluded.

## FLUSH Code Output

Pipe Diameter = 1.77 in.      Flow = 10.00 gpm  
 Total Pipe Length = 106.16 ft      Initial Solid Fraction = 0.500  
 Pipe Roughness = 0.000150      Initial Temperature      25.20 R

Flow, gpm	Flow, lb/s	Final Solid Fraction	Change in x	Pressure Drop, psi
10.000	0.114	0.4859	0.0141	0.0300
15.000	0.171	0.4905	0.0095	0.0655
20.000	0.227	0.4927	0.0073	0.1155
25.000	0.284	0.4939	0.0061	0.1788
30.000	0.341	0.4946	0.0054	0.2558
35.000	0.398	0.4951	0.0049	0.3465
40.000	0.455	0.4953	0.0047	0.4509
45.000	0.512	0.4953	0.0047	0.5690
50.000	0.569	0.4953	0.0047	0.7007
55.000	0.626	0.4951	0.0049	0.8462
60.000	0.682	0.4949	0.0051	1.0053
65.000	0.739	0.4946	0.0054	1.1782
70.000	0.796	0.4943	0.0057	1.3647
75.000	0.853	0.4938	0.0062	1.5649
80.000	0.910	0.4934	0.0066	1.7789
85.000	0.967	0.4928	0.0072	2.0065
90.000	1.024	0.4923	0.0077	2.2479
95.000	1.081	0.4916	0.0084	2.5029
100.000	1.137	0.4910	0.0090	2.7717

TABLE 10. - EXAMPLE FLUSH CODE INPUT FOR HEAT LEAK CALCULATIONS FOR STANDARD INSULATION AND BARE PIPE ON ELEMENTS 3, 6, 12, 15, AND 18

\$PARAMS

FLOWIN=50.0, FLOWMAX=90.0, FLOWINC=10.0, E = 0.00015,  
P = 35., X1 = 0.5,  
ZIN=25., NUM=18, OUT=1.,  
LONG=+1., XPUT=1., FRICT=+1.0, HEAT=+1.0,  
TAMB=540., VAC=2.0, XRS=10., NOJAC=5

\$END

ELE	TYPE	DIA, in.	LENGTH, ft	TEMP., R.	Q	K - ft/D	HT., ft
1	PI	1.770	1.000	25.20	-1.00	0.0	25.0
2	VA	1.770	0.500	25.20	27.30	7.58	25.0
3	PI	1.770	19.00	25.20	0.633	0.0	25.0
4	BE	1.770	1.000	25.20	1.899	0.0	25.0
5	BA	1.770	0.790	25.20	14.78	0.0	25.0
6	PI	1.770	19.00	25.20	0.633	0.0	25.0
7	BE	1.770	1.000	25.20	1.899	0.0	25.0
8	BA	1.770	0.790	25.20	14.78	0.0	25.0
9	PI	1.770	19.00	25.20	-1.00	0.0	25.0
10	BE	1.770	1.000	25.20	1.899	0.0	25.0
11	BA	1.770	0.790	25.20	14.78	0.0	25.0
12	PI	1.770	19.00	25.20	0.633	0.0	25.0
13	BE	1.770	1.000	25.20	1.899	0.0	25.0
14	BA	1.770	0.790	25.20	14.78	0.0	25.0
15	PI	1.770	19.00	25.20	0.633	0.0	25.0
16	BE	1.770	1.000	25.20	1.899	0.0	25.0
17	VA	1.770	0.500	25.20	27.30	7.58	25.0
18	PI	1.770	1.000	25.20	0.633	0.0	25.0

Retention Heat Transfer Parameters

ELE	O.D.	I.D.	Ei	EO
1	1.900	3.334	0.30	0.30
9	1.900	3.334	0.30	0.30

Conduction Heat Transfer Parameters

ELE	O.D.	T1	T2	T3
3	1.900	0.125	0.125	0.125
6	1.900			
12	1.900	0.125	0.125	0.125
15	1.900	0.125	0.125	
18	1.900	0.125		

ELE	PIPE CODE	MAT. 1 CODE	MAT. 2 CODE	MAT. 3 CODE
3	102	201	203	205
6	102			
12	101	202	204	301
15	103	201	203	
18	104	101		

TABLE 11. - MATERIAL CODES FOR  
USE IN STANDARD INSULATION  
OR BARE PIPE CALCULATIONS

Material code	Material
101	Teflon
102	304 Stainless steel
103	Carbon steel
104	6063 Aluminum
201	Foam glass
202	Silica
203	Polyurethane
204	Polystyrene
205	Foam rubber
301	Fiber glass
302	Perlite

TABLE 12. - EXAMPLE FLUSH CODE OUTPUT FOR HEAT LEAK CALCULATIONS FOR  
STANDARD INSULATION AND BARE PIPE ON ELEMENTS 3, 6, 12, 15, AND 18

\*\*\*\*\*

Input Parameters

Element	Type	Dia., in.	Length, ft	Temp., R	Q, Btu/hr-ft	K=f L/D
1	PI	1.770	1.000	25.20	0.552	0.000
2	VA	1.770	0.500	25.20	27.300	7.580
3	PI	1.770	19.000	25.20	32.156	0.000
4	BE	1.770	1.000	25.20	1.899	0.000
5	BA	1.770	0.790	25.20	14.780	0.000
6	PI	1.770	19.000	25.20	253.611	0.000
7	BE	1.770	1.000	25.20	1.899	0.000
8	BA	1.770	0.790	25.20	14.780	0.000
9	PI	1.770	19.000	25.20	0.552	0.000
10	BE	1.770	1.000	25.20	1.899	0.000
11	BA	1.770	0.790	25.20	14.780	0.000
12	PI	1.770	19.000	25.20	35.507	0.000
13	BE	1.770	1.000	25.20	1.899	0.000
14	BA	1.770	0.790	25.20	14.780	0.000
15	PI	1.770	19.000	25.20	37.947	0.000
16	BE	1.770	1.000	25.20	1.899	0.000
17	VA	1.770	0.500	25.20	27.300	7.580
18	PI	1.770	1.000	25.20	208.666	0.000

Radiation Heat Transfer Parameters

Element	Inner O.D.	Outer I.D.	E <sub>i</sub>	E <sub>o</sub>
1	1.900	3.334	0.30000	0.30000
9	1.900	3.334	0.30000	0.30000

TABLE 12. - Concluded.

## Bare Pipe or Standard Insulation Parameters

Element	Pipe O.D.	Insulation 1 Thickness, ft	Insulation 2 Thickness, ft	Insulation 3 Thickness, ft
3	1.9000	0.1250	0.1250	0.1250
6	1.9000	0.0000	0.0000	0.0000
12	1.9000	0.1250	0.1250	0.1250
15	1.9000	0.1250	0.1250	0.0000
18	1.9000	0.1250	0.0000	0.0000

Element	Pipe Material	Insulation 1	Insulation 2	Insulation 3
3	304 St Steel	Foam Glass	Polyurethane	Foam Rubber
6	304 St Steel			
12	Teflon	Silica	Polystyrene	Fiber Glass
15	Carbon Steel	Foam Glass	Polyurethane	
18	6063 Alum.	Teflon		

\*\*\*\*\*  
FLUSH Code Output

.....  
 Pipe Diameter = 1.77 in.      Flow = 50.00 gpm  
 Total Pipe Length = 106.16 ft      Initial Solid Fraction = 0.500  
 Pipe Roughness = 0.000150      Initial Temperature = 25.20 R  
 .....

Flow, gpm	Flow, lb/s	Final Solid Fraction	Change in x	Pressure Drop, psi
50.000	0.569	0.3589	0.1411	0.7080
60.000	0.682	0.3761	0.1239	1.0145
70.000	0.796	0.3884	0.1116	1.3758
80.000	0.910	0.3977	0.1023	1.7920
90.000	1.024	0.4049	0.0951	2.2630



TABLE 13. - EXAMPLE FLUSH CODE INPUT WITH PUMPS AT ELEMENTS 2 AND 4

\$PARAMS

FLOWIN=352.3, FLOWMAX=352.3, FLOWINC=10.0, E = 0.00015,

P = 25., X1 = 0.5,

ZIN=25., NUM=5, OUT=1.,

LONG=+1., XPUT=1., FRICT=+1.0, HEAT=+1.0,

TAMB=540., VAC=0.0, XRS=10., NOJAC=0

\$END

ELE	TYPE	DIA, in.	LENGTH, ft	TEMP., R.	Q	K=fL/D	HT., ft
1	PI	1.770	5.000	25.20	0.633	0.0	25.0
2	PH	1.770	1.000	25.20	1.690	.322	25.0
3	PI	1.770	5.000	25.20	0.633	0.0	25.0
4	PP	1.770	1.000	25.20	2.643	.322	25.0
5	PI	1.770	5.000	25.20	0.633	0.0	25.0

TABLE 14. - EXAMPLE FLUSH CODE OUTPUT WITH PUMPS AT ELEMENTS 2 AND 4

\*\*\*\*\*

Input Parameters

Element	Type	Dia., in.	Length, ft	Temp., R	Q, Btu/hr-ft.	K=f L/D
1	PI	1.770	5.000	25.20	0.633	0.000
2	PH	1.770	1.000	25.20	1.690	0.322
3	PI	1.770	5.000	25.20	0.633	0.000
4	PP	1.770	1.000	25.20	2.643	0.322
5	PI	1.770	5.000	25.20	0.633	0.000

FLUSH1 Code Output

Pipe Diameter = 1.77 in.      Flow = 352.30 gpm  
 Total Pipe Length = 17.00 ft      Initial Solid Fraction = 0.500  
 Pipe Roughness = 0.000150      Initial Temperature = 25.20 R

Node	Pressure Drop, psi	Flow, lb/s	Final Solid Fraction	Change in x	Pressure, psi
1	0.00000	4.0056	0.5000	0.0000	25.0000
2	0.78304	4.0056	0.4978	0.0022	24.2170
3	-1.86134	4.0056	0.4893	0.0107	26.8613
4	-1.07737	4.0056	0.4872	0.0128	26.0774
5	-3.72037	4.0056	0.4787	0.0213	28.7204
6	-2.93547	4.0056	0.4765	0.0235	27.9355

TABLE 15. - CRITICAL  
FLOW RATES FOR SLUSH  
PIPING

[Based on a critical velocity of 1.5 ft/sec;  
Data in the table are  
scaled from experimental  
results obtained  
using a 0.652-in.-i.d.  
pipe (ref. 2).]

Nominal diameter, in.	Critical flow rate, gpm
1	5.2
1.5	11.5
2	18.5
2.5	27.9
3	40.8
4	69.0
5	104.9
6	150.7
8	259.5
10	403.4
12	568.0

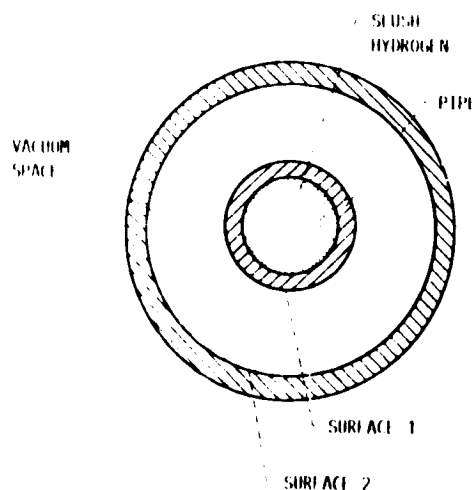


FIGURE 1. CROSS SECTION OF VACUUM JACKETED PIPE.

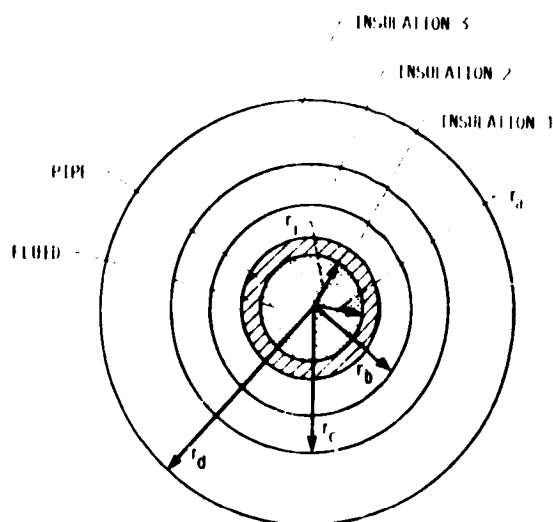


FIGURE 2. CROSS SECTION OF PIPE WITH THREE LAYERS OF INSULATION.

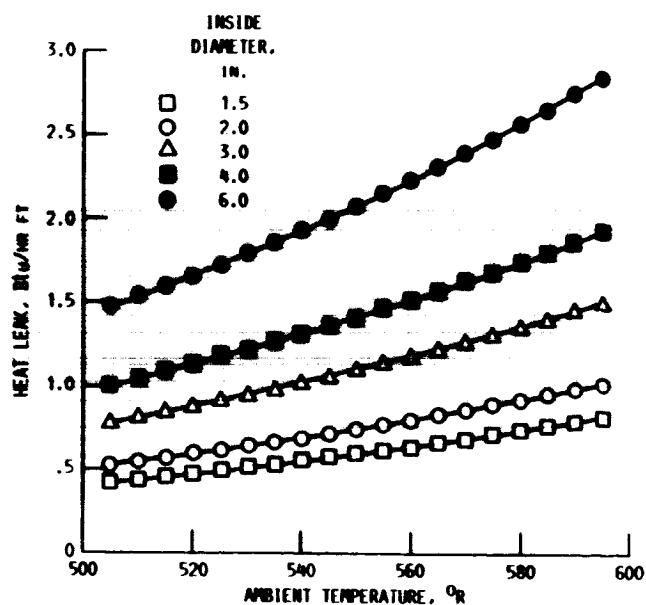


FIGURE 3. - HEAT LEAK VERSUS AMBIENT TEMPERATURE FOR SEVERAL SIZES OF VACUUM JACKETED PIPE. NUMBER OF RADIATION SHIELDS, 10; EMISSIVITY, 0.3.

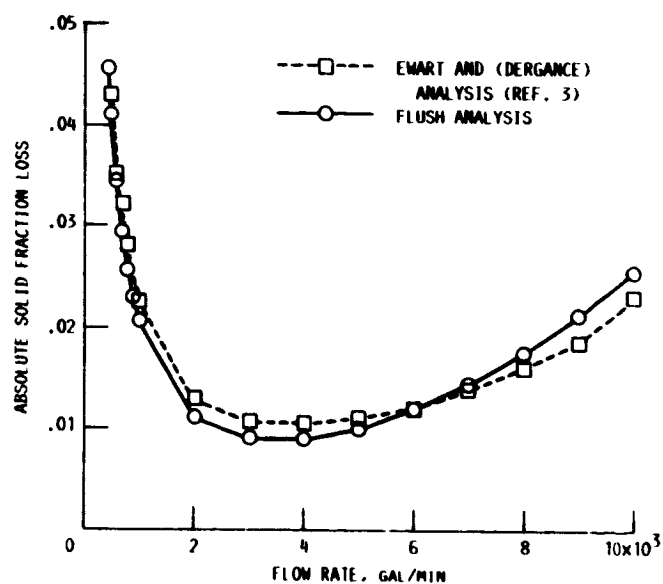
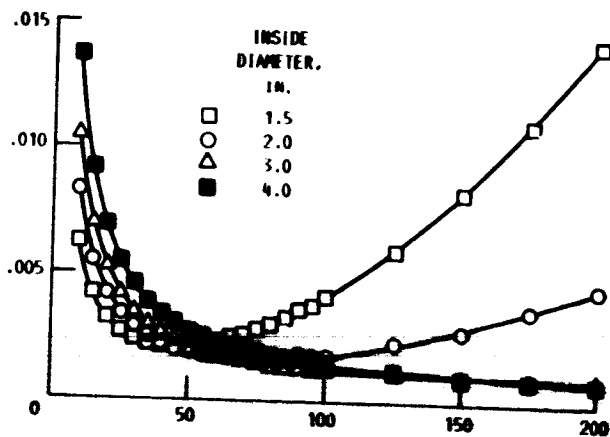
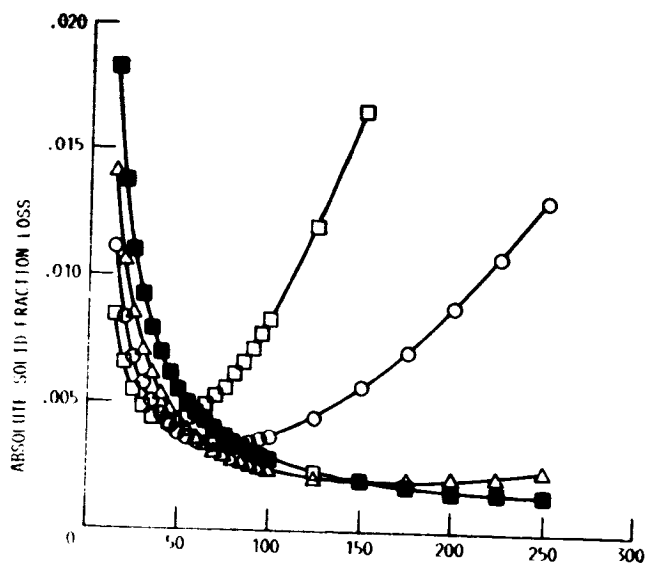


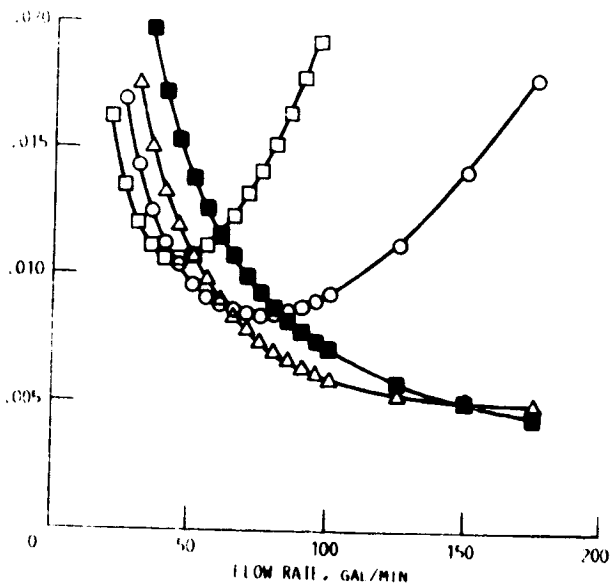
FIGURE 4. - COMPARISON OF FLUSH ANALYSIS WITH PREVIOUS SLUSH HYDROGEN ANALYSIS. PIPE 12-IN.-DIAMETER SCHEDULE 5S; HEAT LEAK INTO ELEMENT,  $Q$ , 12 Btu/hr ft; PIPE LENGTH, 1750 FT.



(a) TOTAL PIPE LENGTH, 100 FT.



(b) TOTAL PIPE LENGTH, 200 FT.



(c) TOTAL PIPE LENGTH, 500 FT.

FIGURE 5. SLUSH LOSS VERSUS FLOW RATE FOR SEVERAL SIZES OF VACUUM JACKETED PIPE.

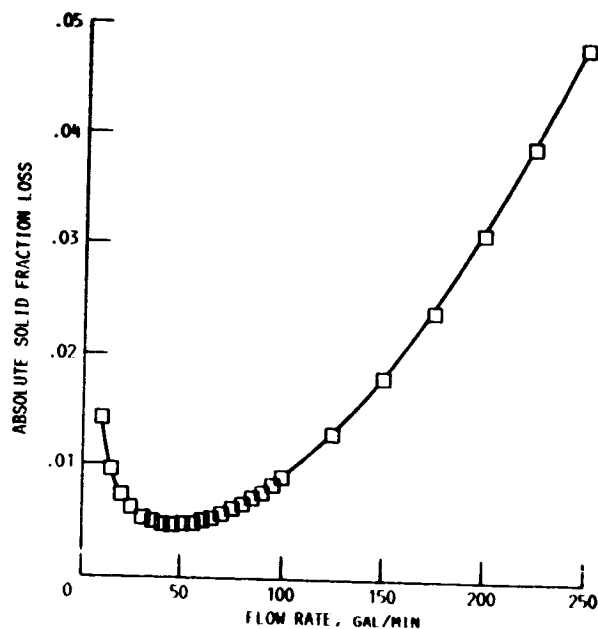


FIGURE 6. - EXPECTED SLUSH SOLID FRACTION LOSS AT K-SITE FOR 1.5-IN. DIAMETER VACUUM-JACKETED PIPE.

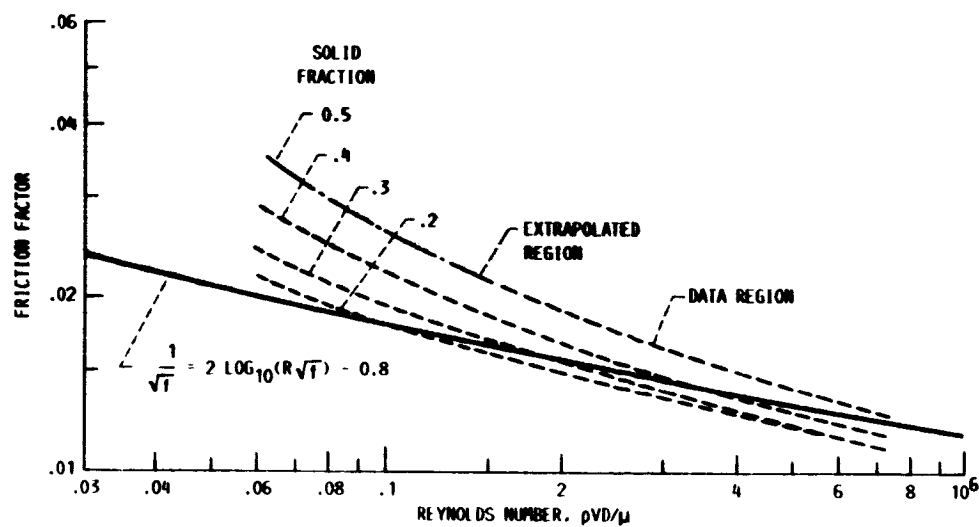


FIGURE 7. - FRICTION FACTOR VERSUS REYNOLDS NUMBER FOR SLUSH HYDROGEN. PIPE DIAMETER, D, 0.652 IN.; TRIPLE-POINT HYDROGEN VISCOSITY,  $\mu = 1.73 \times 10^{-5}$  LBW/FT SEC. (FROM REF. 2.)

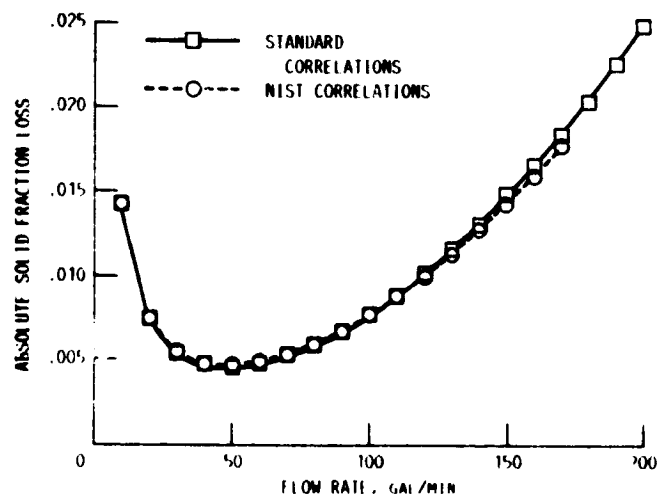


FIGURE 8. COMPARISON OF SLUSH HYDROGEN LOSS WITH VARIOUS FRICTION FACTOR CORRELATIONS. 1.5 IN. I.D. VACUUM JACK ETTED PIPE; INITIAL SLUSH SOLID FRACTION, 0.5.



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